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NEWS 1	Web Page URLs for STN Seminar Schedule - N. America
NEWS 2	"Ask CAS" for self-help around the clock
NEWS 3	May 12 EXTEND option available in structure searching
NEWS 4	May 12 Polymer links for the POLYLINK command completed in REGISTRY
NEWS 5	May 27 New UPM (Update Code Maximum) field for more efficient patent SDIs in CAplus
NEWS 6	May 27 CAplus super roles and document types searchable in REGISTRY
NEWS 7	Jun 28 Additional enzyme-catalyzed reactions added to CASREACT
NEWS 8	Jun 28 ANTE, AQUALINE, BIOENG, CIVILENG, ENVIROENG, MECHENG, and WATER from CSA now available on STN(R)
NEWS 9	Jul 12 BEILSTEIN enhanced with new display and select options, resulting in a closer connection to BABS
NEWS 10	Jul 30 BEILSTEIN on STN workshop to be held August 24 in conjunction with the 228th ACS National Meeting
NEWS 11	AUG 02 IFIPAT/IFIUDB/IFICDB reloaded with new search and display fields
NEWS 12	AUG 02 CAplus and CA patent records enhanced with European and Japan Patent Office Classifications
NEWS 13	AUG 02 STN User Update to be held August 22 in conjunction with the 228th ACS National Meeting
NEWS 14	AUG 02 The Analysis Edition of STN Express with Discover! (Version 7.01 for Windows) now available
NEWS 15	AUG 04 Pricing for the Save Answers for SciFinder Wizard within STN Express with Discover! will change September 1, 2004
NEWS EXPRESS	JULY 30 CURRENT WINDOWS VERSION IS V7.01, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 11 AUGUST 2004
NEWS HOURS	STN Operating Hours Plus Help Desk Availability
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FILE 'HOME' ENTERED AT 10:54:23 ON 18 AUG 2004

=> file reg  
 COST IN U.S. DOLLARS  
 FULL ESTIMATED COST

SINCE FILE ENTRY	TOTAL SESSION
0.21	0.21

FILE 'REGISTRY' ENTERED AT 10:54:32 ON 18 AUG 2004  
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STRUCTURE FILE UPDATES: 17 AUG 2004 HIGHEST RN 727974-89-2  
 DICTIONARY FILE UPDATES: 17 AUG 2004 HIGHEST RN 727974-89-2

TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

Please note that search-term pricing does apply when  
 conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

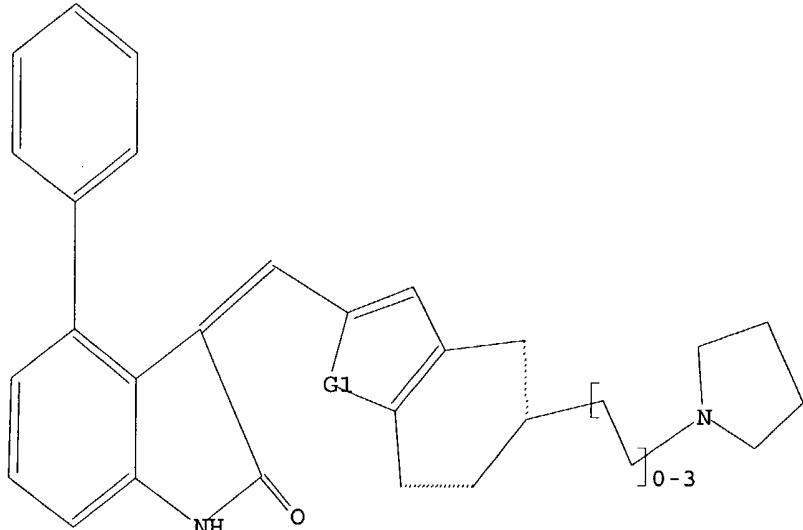
Experimental and calculated property data are now available. For more  
 information enter HELP PROP at an arrow prompt in the file or refer  
 to the file summary sheet on the web at:

<http://www.cas.org/ONLINE/DBSS/registryss.html>

=>  
 Uploading c:\program files\stnexp\queries\10736243.1

L1 STRUCTURE UPLOADED

=> d 11  
 L1 HAS NO ANSWERS  
 L1 STR



G1 CH,NH

Structure attributes must be viewed using STN Express query preparation.

=> s 11 sss full  
 FULL SEARCH INITIATED 10:55:00 FILE 'REGISTRY'  
 FULL SCREEN SEARCH COMPLETED - 109 TO ITERATE

100.0% PROCESSED 109 ITERATIONS 2 ANSWERS  
 SEARCH TIME: 00.00.01

L2 2 SEA SSS FUL L1

=> file caplus	SINCE FILE	TOTAL
COST IN U.S. DOLLARS	ENTRY	SESSION
FULL ESTIMATED COST	155.42	155.63

FILE 'CAPLUS' ENTERED AT 10:55:05 ON 18 AUG 2004  
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FILE COVERS 1907 - 18 Aug 2004 VOL 141 ISS 8  
 FILE LAST UPDATED: 17 Aug 2004 (20040817/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s 12  
 L3 1 L2

=> d 13 fbib hitstr abs total

L3 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2004 ACS on STN  
 AN 2002:539677 CAPLUS  
 DN 137:109202  
 TI Preparation of 4-aryl substituted indolinones as protein kinase signal transduction modulators for inhibiting abnormal cell proliferation  
 IN Cui, Jingrong; Zhang, Ruofei; Shen, Hong; Chu, Ji Yu; Zhang, Fang-Jie; Koenig, Marcel; Do, Steven Huy; Li, Xiaoyuan; Wei, Chung Chen; Tang, Peng Cho  
 PA USA  
 SO PCT Int. Appl., 560 pp.  
 CODEN: PIXXD2  
 DT Patent  
 LA English

## FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2002055517	A2	20020718	WO 2001-US48564	20011220
	WO 2002055517	A3	20020926		
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	RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			US 2000-256479P	P 20001220
	US 2003069297	A1	20030410	US 2001-23488	20011220
	US 6677368	B2	20040113	US 2000-256479P	P 20001220
	EP 1349852	A2	20031008	EP 2001-997065	20011220
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR			US 2000-256479P	P 20001220
				WO 2001-US48564	W 20011220
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				US 2000-256479P	P 20001220
				WO 2001-US48564	W 20011220
	US 2004157909	A1	20040812	US 2003-736243	20031216
				US 2000-256479P	P 20001220
				US 2001-23488	A3 20011220

OS MARPAT 137:109202

IT 442559-47-9P 442559-48-0P

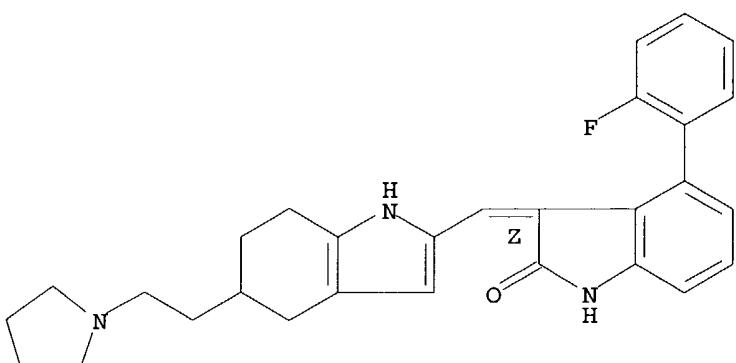
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(target compound; preparation of (aryl)(pyrrolylmethylene)indolinones as protein kinase signal transduction modulators)

RN 442559-47-9 CAPLUS

CN 2H-Indol-2-one, 4-(2-fluorophenyl)-1,3-dihydro-3-[[4,5,6,7-tetrahydro-5-[2-(1-pyrrolidinyl)ethyl]-1H-indol-2-yl]methylene]-, (3Z)- (9CI) (CA INDEX NAME)

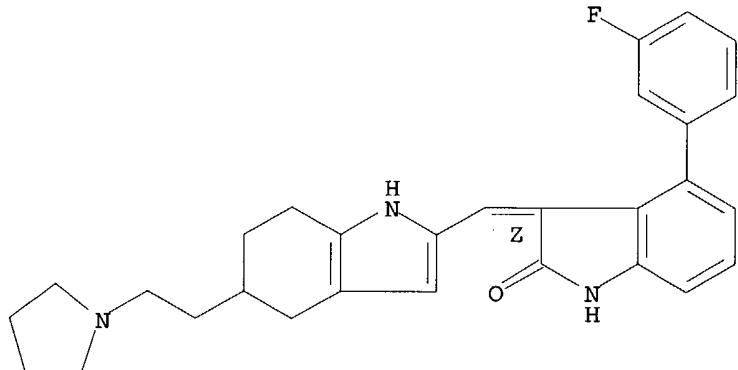
Double bond geometry as shown.



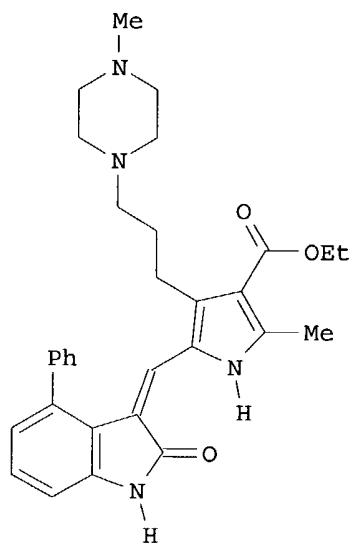
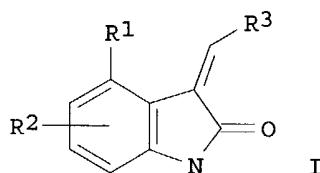
RN 442559-48-0 CAPLUS

CN 2H-Indol-2-one, 4-[(3-fluorophenyl)-1,3-dihydro-3-[(4,5,6,7-tetrahydro-5-[2-(1-pyrrolidinyl)ethyl]-1H-indol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



GI



AB Title compds. I [R1 = (un)substituted aryl or heteroaryl; R2 = H, halo, alkyl, alkenyl, alkynyl, heterocyclyl, etc.; R3 = (un)substituted pyrrole or cycloalkenylpyrrole], as well as pharmaceutical compns. thereof, are prepared and disclosed as compds. capable of modulating protein kinase signal transduction in order to regulate, modulate and/or inhibit abnormal cell proliferation. Thus II, was prepared via condensation of 4-phenyl-1,3-dihydroindol-2-one with 5-formyl-2-methyl-4-[3-(4-methylpiperazin-1-yl)propyl]-1H-pyrrole-3-carboxylic acid Et ester. I were evaluated against eight specific kinases, e.g., FGFR1, for which I possessed IC50 values ( $\mu$ M) of 0.0091-2.07. The present invention also relates to methods for treating protein kinase related disorders.

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PASSWORD :

TERMINAL (ENTER 1, 2, 3, OR ?):2

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NEWS	4	May 12 Polymer links for the POLYLINK command completed in REGISTRY
NEWS	5	May 27 New UPM (Update Code Maximum) field for more efficient patent SDIs in CAplus
NEWS	6	May 27 CAplus super roles and document types searchable in REGISTRY
NEWS	7	Jun 28 Additional enzyme-catalyzed reactions added to CASREACT
NEWS	8	Jun 28 ANTE, AQUALINE, BIOENG, CIVILENG, ENVIROENG, MECHENG, and WATER from CSA now available on STN(R)
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NEWS	10	Jul 30 BEILSTEIN on STN workshop to be held August 24 in conjunction with the 228th ACS National Meeting
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NEWS	15	AUG 04 Pricing for the Save Answers for SciFinder Wizard within STN Express with Discover! will change September 1, 2004
NEWS EXPRESS		JULY 30 CURRENT WINDOWS VERSION IS V7.01, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 11 AUGUST 2004
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NEWS LOGIN		Welcome Banner and News Items
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=> fdile reg  
FDILE IS NOT A RECOGNIZED COMMAND  
The previous command name entered was not recognized by the system.  
For a list of commands available to you in the current file, enter  
"HELP COMMANDS" at an arrow prompt (=>).

=>  
Uploading  
THIS COMMAND NOT AVAILABLE IN THE CURRENT FILE  
Do you want to switch to the Registry File?  
Choice (Y/n) :  
Switching to the Registry File...  
Some commands only work in certain files. For example, the EXPAND  
command can only be used to look at the index in a file which has an  
index. Enter "HELP COMMANDS" at an arrow prompt (=>) for a list of  
commands which can be used in this file.

=> FILE REGISTRY

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.42	0.42

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STRUCTURE FILE UPDATES: 17 AUG 2004 HIGHEST RN 727974-89-2  
DICTIONARY FILE UPDATES: 17 AUG 2004 HIGHEST RN 727974-89-2

TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

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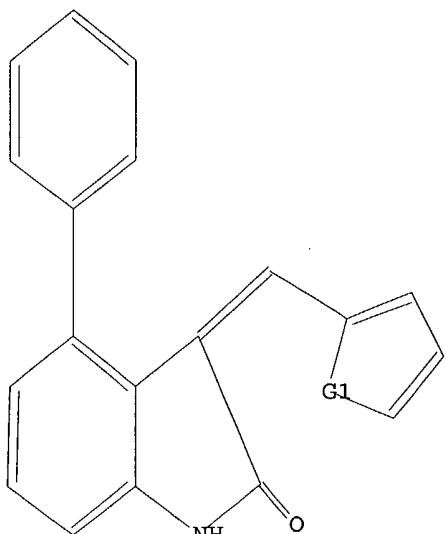
Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more  
information enter HELP PROP at an arrow prompt in the file or refer  
to the file summary sheet on the web at:  
<http://www.cas.org/ONLINE/DBSS/registryss.html>

=>  
Uploading c:\program files\stnexp\queries\10736243.2

L1 STRUCTURE UPLOADED

=> d 11  
L1 HAS NO ANSWERS  
L1 STR



G1 CH,NH

Structure attributes must be viewed using STN Express query preparation.

=> s 11 sss full  
 FULL SEARCH INITIATED 11:18:33 FILE 'REGISTRY'  
 FULL SCREEN SEARCH COMPLETED - 3216 TO ITERATE

100.0% PROCESSED 3216 ITERATIONS 498 ANSWERS  
 SEARCH TIME: 00.00.01

L2 498 SEA SSS FUL L1

=> file caplus			
COST IN U.S. DOLLARS	SINCE FILE	TOTAL	
FULL ESTIMATED COST	ENTRY	SESSION	
	155.42	155.84	

FILE 'CAPLUS' ENTERED AT 11:18:39 ON 18 AUG 2004  
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FILE COVERS 1907 - 18 Aug 2004 VOL 141 ISS 8  
 FILE LAST UPDATED: 17 Aug 2004 (20040817/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s 12  
L3 3 L2

=> d 13 fbib hitstr abs total

L3 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2004 ACS on STN  
AN 2004:182368 CAPLUS  
DN 140:229401  
TI Three hybrid assay system for isolating ligand-binding polypeptides and for isolating small mol. ligands  
IN Come, Jon H.; Becker, Frank; Kley, Nikolai A.; Reichel, Christoph  
PA USA  
SO U.S. Pat. Appl. Publ., 238 pp., Cont.-in-part of U.S. Ser. No. 91,177.  
CODEN: USXXCO

DT Patent  
LA English  
FAN.CNT 3

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	US 2004043388	A1	20040304	US 2002-234985 US 2001-272932P US 2001-278233P US 2001-329437P US 2002-91177	20020903 P 20010302 P 20010323 P 20011015 A2 20020304
	US 2003165873	A1	20030904	US 2002-91177 US 2001-272932P US 2001-278233P US 2001-329437P	20020304 P 20010302 P 20010323 P 20011015

PATENT FAMILY INFORMATION:

FAN 2002:696096

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2002070662	A2	20020912	WO 2002-US6677	20020304
	WO 2002070662	A3	20021227		
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	RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			US 2001-272932P US 2001-278233P US 2001-329437P	P 20010302 P 20010323 P 20011015
	EP 1364212	A2	20031126	EP 2002-723332	20020304
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FAN 2003:319902

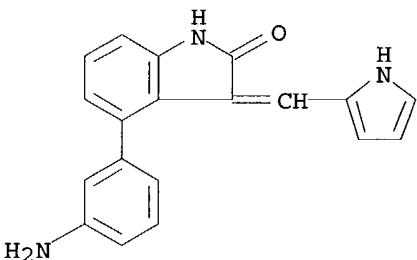
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	WO 2003033499	A3	20030814		
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	RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			US 2001-329437P	P 20011015
				US 2001-336962P	P 20011203
EP 1446405	A2	20040818	EP 2002-797047	20021015	
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				US 2001-336962P	P 20011203
				WO 2002-US33052	W 20021015
US 2003162797	A1	20030828	US 2002-321284	20021217	
US 6753329	B2	20040622		US 2001-336962P	P 20011203
				WO 2002-US33052	A1 20021015

IT 666838-05-7D, conjugates

RL: BUU (Biological use, unclassified); THU (Therapeutic use); BIOL (Biological study); USES (Uses)  
 (three hybrid assay system for isolating ligand-binding polypeptides and for isolating small mol. ligands)

RN 666838-05-7 CAPLUS

CN 2H-Indol-2-one, 4-(3-aminophenyl)-1,3-dihydro-3-(1H-pyrrol-2-ylmethylene)-(9CI) (CA INDEX NAME)



AB The invention provides compns. and methods for isolating ligand-binding polypeptides for a user-specified ligand, and for isolating small mol. ligands for a user-specified target polypeptide using an improved class of hybrid ligand compds. Preparation of compds., e.g. a methotrexate moiety linked by a polyethylene glycol moiety to dexamethasone, is described.

L3 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2004 ACS on STN

AN 2002:539677 CAPLUS

DN 137:109202

TI Preparation of 4-aryl substituted indolinones as protein kinase signal

IN transduction modulators for inhibiting abnormal cell proliferation  
 Cui, Jingrong; Zhang, Ruofei; Shen, Hong; Chu, Ji Yu; Zhang, Fang-Jie;  
 Koenig, Marcel; Do, Steven Huy; Li, Xiaoyuan; Wei, Chung Chen; Tang, Peng  
 Cho

PA USA

SO PCT Int. Appl., 560 pp.

CODEN: PIXXD2

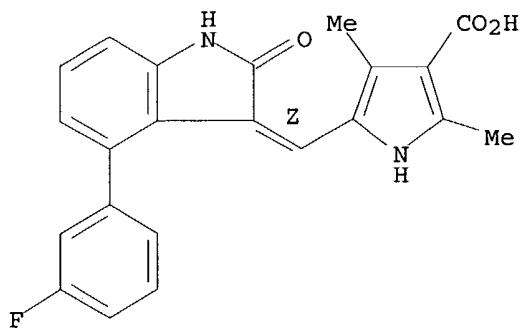
DT Patent

LA English

FAN.CNT 1

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI WO 2002055517	A2	20020718	WO 2001-US48564	20011220
WO 2002055517	A3	20020926		
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
			US 2000-256479P	P 20001220
US 2003069297	A1	20030410	US 2001-23488	20011220
US 6677368	B2	20040113	US 2000-256479P	P 20001220
EP 1349852	A2	20031008	EP 2001-997065	20011220
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR				
US 2000-256479P P 20001220				
WO 2001-US48564 W 20011220				
JP 2004518669	T2	20040624	JP 2002-556186	20011220
US 2000-256479P P 20001220				
WO 2001-US48564 W 20011220				
US 2004157909	A1	20040812	US 2003-736243	20031216
US 2000-256479P P 20001220				
US 2001-23488 A3 20011220				
OS MARPAT 137:109202				
IT 442563-63-5				
RL: RCT (Reactant); RACT (Reactant or reagent) (preparation of (aryl)(pyrrolylmethylene)indolinones as protein kinase signal transduction modulators)				
RN 442563-63-5	CAPLUS			
CN 1H-Pyrrole-3-carboxylic acid, 5-[(Z)-[4-(3-fluorophenyl)-1,2-dihydro-2-oxo- 3H-indol-3-ylidene]methyl]-2,4-dimethyl- (9CI) (CA INDEX NAME)				

Double bond geometry as shown.



IT	442558-03-4P	442558-04-5P	442558-05-6P
	442558-06-7P	442558-07-8P	442558-08-9P
	442558-09-0P	442558-10-3P	442558-11-4P
	442558-12-5P	442558-13-6P	442558-14-7P
	442558-15-8P	442558-16-9P	442558-17-0P
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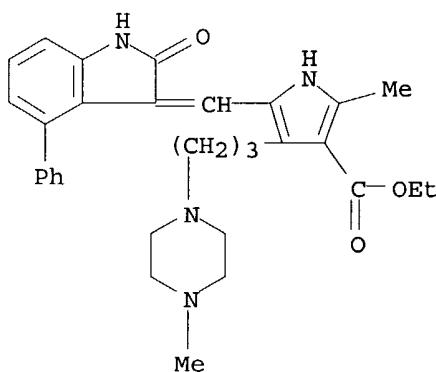
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RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(target compound; preparation of (aryl)(pyrrolylmethylene)indolinones as protein kinase signal transduction modulators)

RN 442558-03-4 CAPLUS

CN 1H-Pyrrole-3-carboxylic acid, 5-[(1,2-dihydro-2-oxo-4-phenyl-3H-indol-3-ylidene)methyl]-2-methyl-4-[3-(4-methyl-1-piperazinyl)propyl]-, ethyl ester (9CI) (CA INDEX NAME)



AN 2000:421132 CAPLUS

DN 133:43433

TI Preparation of 4-aryl-3-(azolylmethylidene)-2-oxindoles as inhibitors of JNK protein kinases.

IN Corbett, Wendy Lea; Luk, Kin-chun; Mahaney, Paige E.

PA F. Hoffmann-La Roche A.-G., Switz.

SO PCT Int. Appl., 91 pp.

CODEN: PIXXD2

DT Patent

LA English

FAN.CNT 1

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI WO 2000035909	A1	20000622	WO 1999-EP9673	19991209
W: AE, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
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Patel

<8/18/2004>

DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

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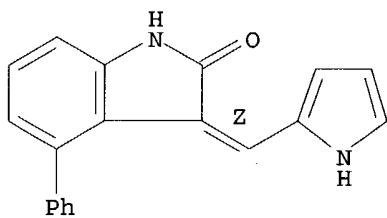
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RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)  
 (preparation of 4-aryl-3-(azolylmethylidene)-2-oxindoles as inhibitors of JNK protein kinases)

RN 276250-95-4 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-phenyl-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

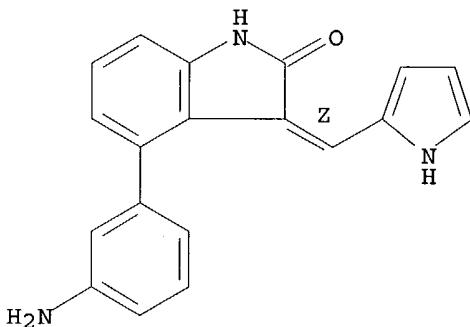
Double bond geometry as shown.



RN 276250-97-6 CAPLUS

CN 2H-Indol-2-one, 4-(3-aminophenyl)-1,3-dihydro-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

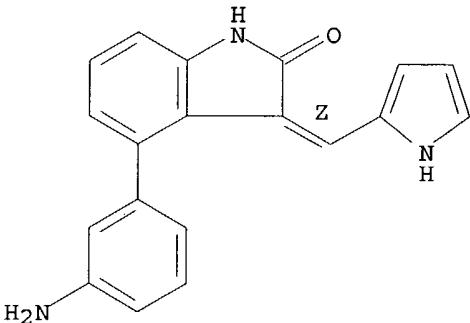
Double bond geometry as shown.



RN 276250-98-7 CAPLUS

CN 2H-Indol-2-one, 4-(3-aminophenyl)-1,3-dihydro-3-(1H-pyrrol-2-ylmethylene)-, monohydrochloride, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

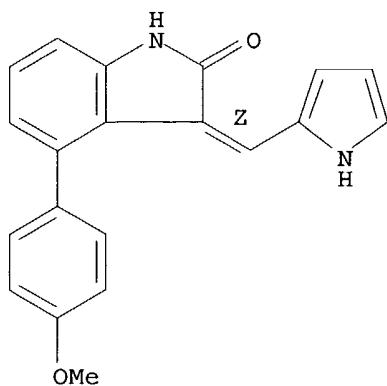


● HCl

RN 276250-99-8 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(4-methoxyphenyl)-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

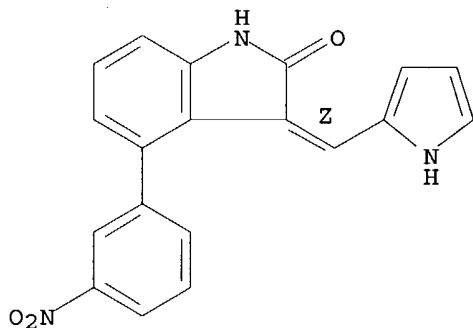
Double bond geometry as shown.



RN 276251-00-4 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(3-nitrophenyl)-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

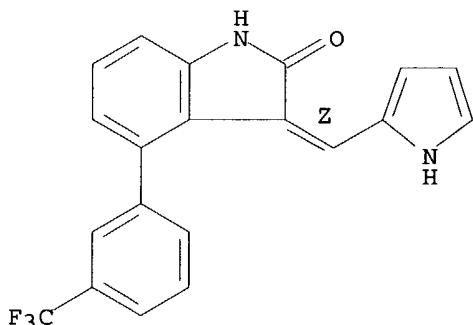
Double bond geometry as shown.



RN 276251-02-6 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-3-(1H-pyrrol-2-ylmethylene)-4-[3-(trifluoromethyl)phenyl]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

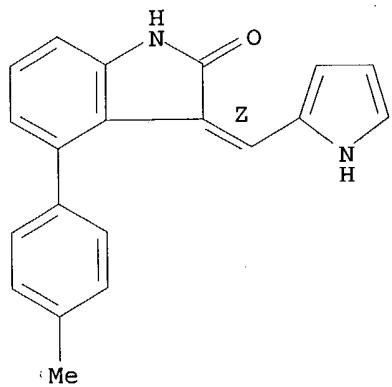


RN 276251-04-8 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(4-methylphenyl)-3-(1H-pyrrol-2-ylmethylene)-

, (3Z) - (9CI) (CA INDEX NAME)

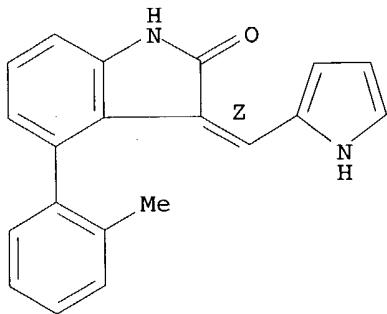
Double bond geometry as shown.



RN 276251-06-0 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(2-methylphenyl)-3-(1H-pyrrol-2-ylmethylene)-, (3Z) - (9CI) (CA INDEX NAME)

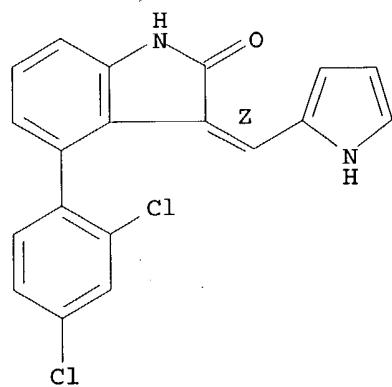
Double bond geometry as shown.



RN 276251-08-2 CAPLUS

CN 2H-Indol-2-one, 4-(2,4-dichlorophenyl)-1,3-dihydro-3-(1H-pyrrol-2-ylmethylene)-, (3Z) - (9CI) (CA INDEX NAME)

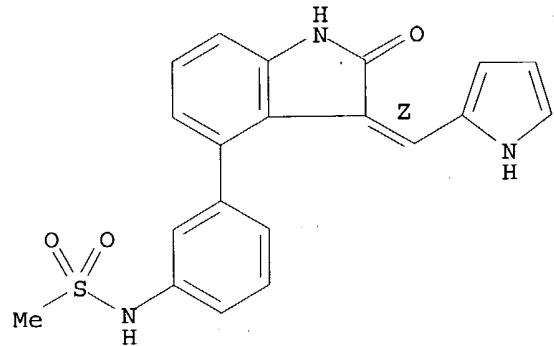
Double bond geometry as shown.



RN 276251-10-6 CAPLUS

CN Methanesulfonamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]- (9CI) (CA INDEX NAME)

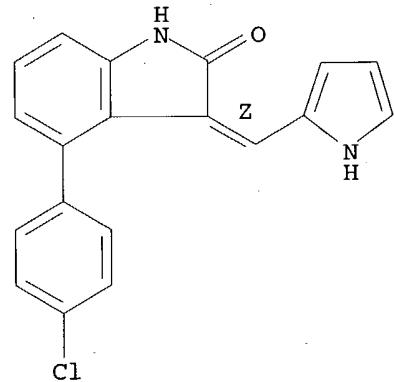
Double bond geometry as shown.



RN 276251-12-8 CAPLUS

CN 2H-Indol-2-one, 4-(4-chlorophenyl)-1,3-dihydro-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

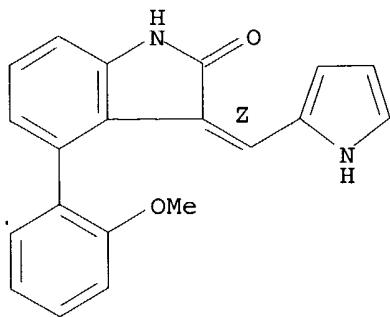
Double bond geometry as shown.



RN 276251-14-0 CAPLUS

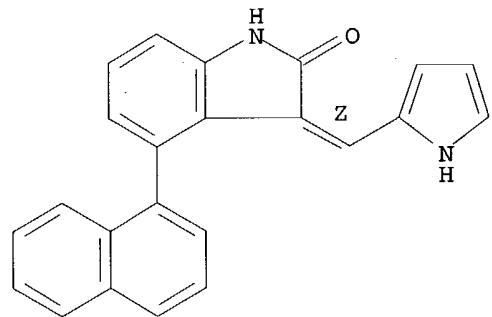
CN 2H-Indol-2-one, 1,3-dihydro-4-(2-methoxyphenyl)-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



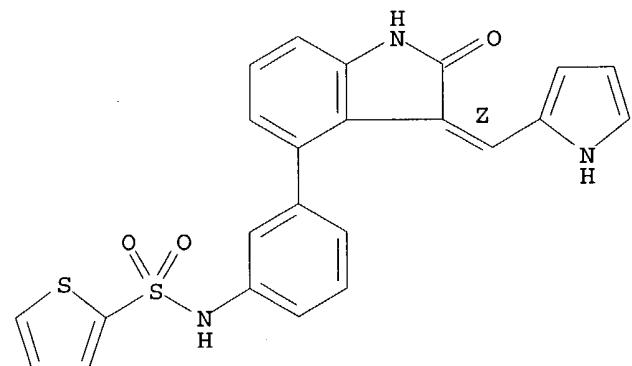
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 CN 2H-Indol-2-one, 1,3-dihydro-4-(1-naphthalenyl)-3-(1H-pyrrol-2-ylmethylene)-  
     , (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



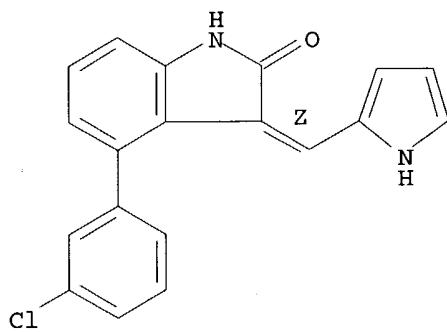
RN 276251-18-4 CAPLUS  
 CN 2-Thiophenesulfonamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-  
     ylmethylene)-1H-indol-4-yl]phenyl]- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 276251-19-5 CAPLUS  
 CN 2H-Indol-2-one, 4-(3-chlorophenyl)-1,3-dihydro-3-(1H-pyrrol-2-ylmethylene)-  
     , (3Z)- (9CI) (CA INDEX NAME)

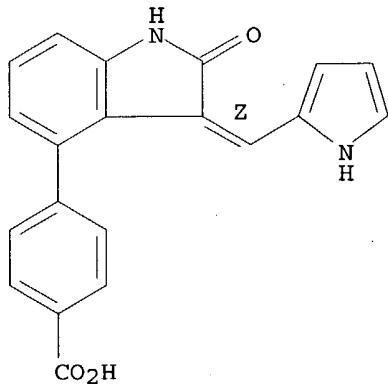
Double bond geometry as shown.



RN 276251-20-8 CAPLUS

CN Benzoic acid, 4-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]- (9CI) (CA INDEX NAME)

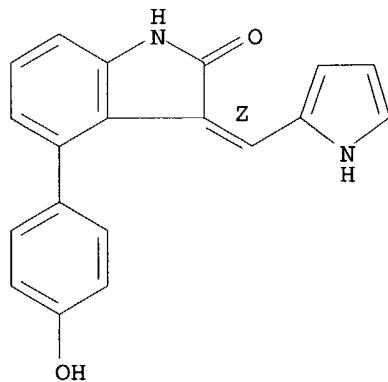
Double bond geometry as shown.



RN 276251-21-9 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(4-hydroxyphenyl)-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

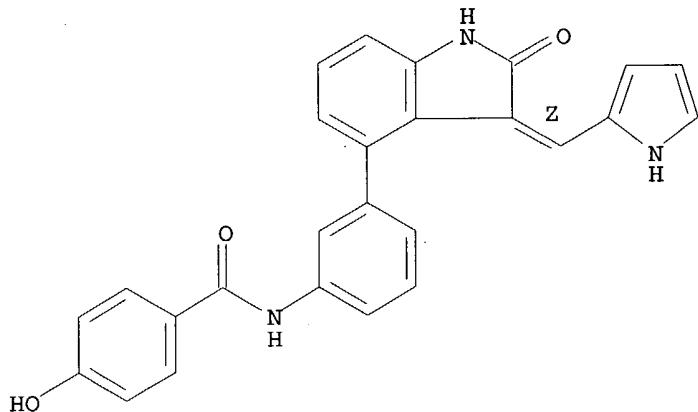
Double bond geometry as shown.



RN 276251-22-0 CAPLUS

CN Benzamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-4-hydroxy- (9CI) (CA INDEX NAME)

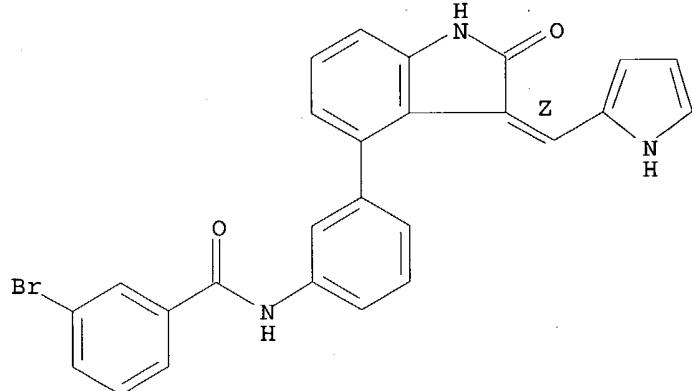
Double bond geometry as shown.



RN 276251-23-1 CAPLUS

CN Benzamide, 3-bromo-N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]- (9CI) (CA INDEX NAME)

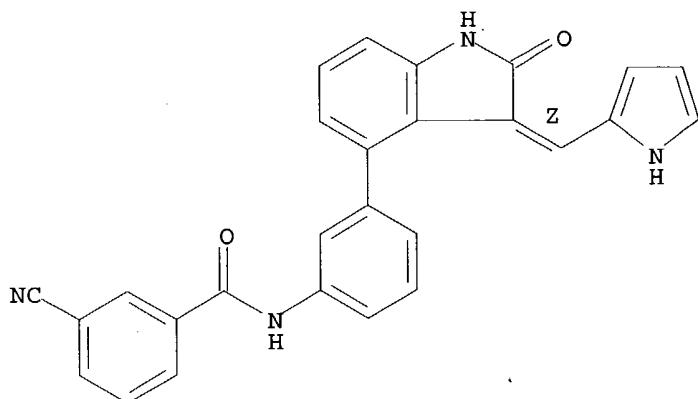
Double bond geometry as shown.



RN 276251-24-2 CAPLUS

CN Benzamide, 3-cyano-N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]- (9CI) (CA INDEX NAME)

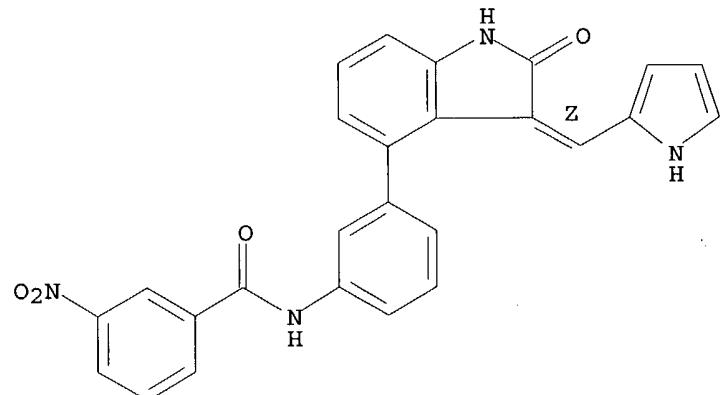
Double bond geometry as shown.



RN 276251-25-3 CAPLUS

CN Benzamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-3-nitro- (9CI) (CA INDEX NAME)

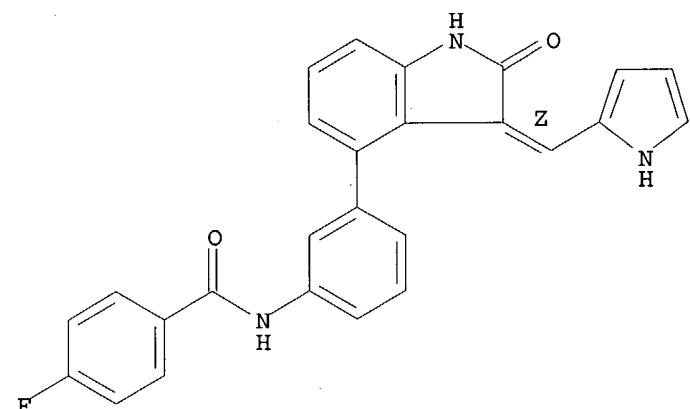
Double bond geometry as shown.



RN 276251-26-4 CAPLUS

CN Benzamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-4-fluoro- (9CI) (CA INDEX NAME)

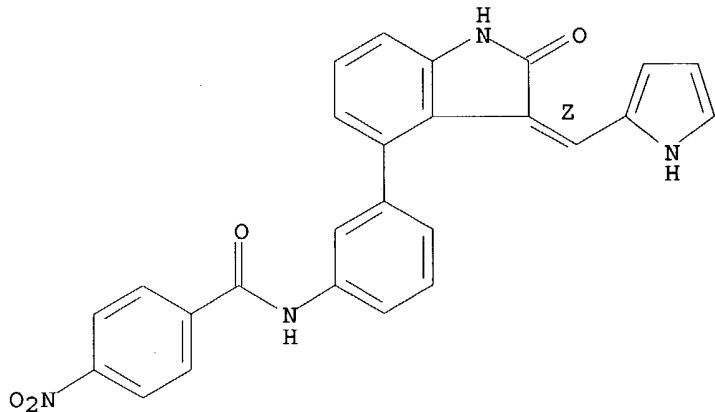
Double bond geometry as shown.



RN 276251-27-5 CAPLUS

CN Benzamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-4-nitro- (9CI) (CA INDEX NAME)

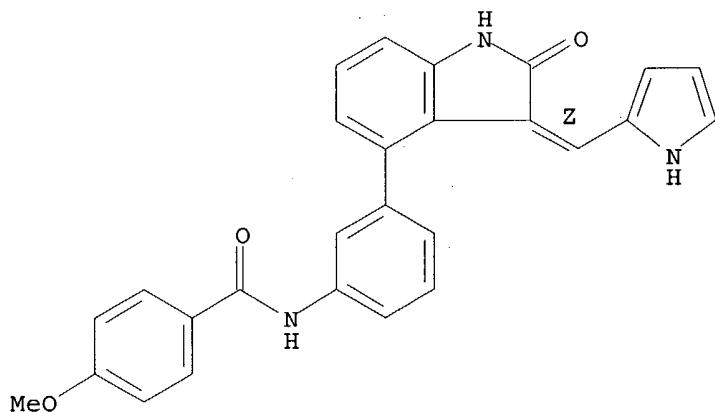
Double bond geometry as shown.



RN 276251-28-6 CAPLUS

CN Benzamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-4-methoxy- (9CI) (CA INDEX NAME)

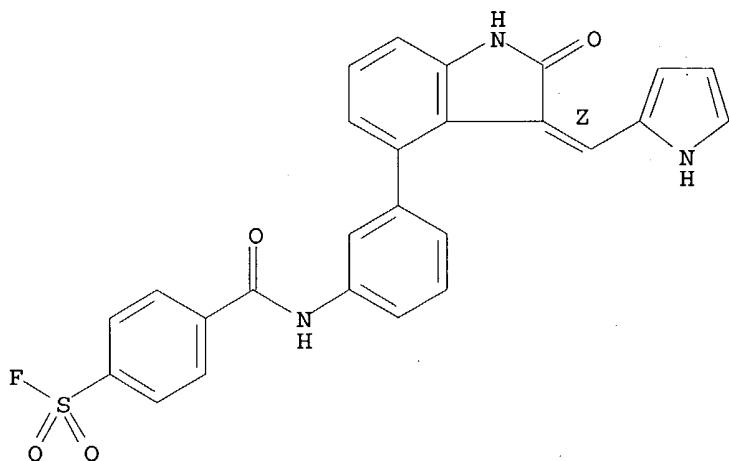
Double bond geometry as shown.



RN 276251-29-7 CAPLUS

CN Benzenesulfonyl fluoride, 4-[[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]amino]carbonyl- (9CI) (CA INDEX NAME)

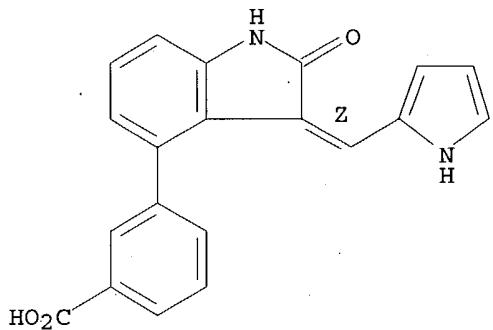
Double bond geometry as shown.



RN 276251-30-0 CAPLUS

CN Benzoic acid, 3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl] - (9CI) (CA INDEX NAME)

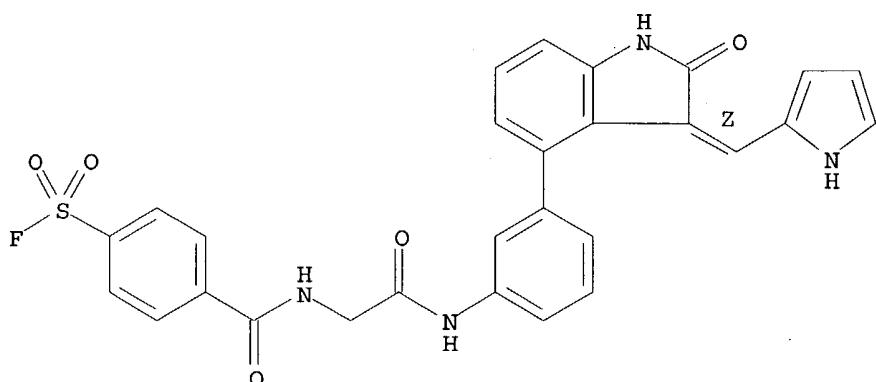
Double bond geometry as shown.



RN 276251-31-1 CAPLUS

CN Benzenesulfonyl fluoride, 4-[[2-[[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]amino]-2-oxoethyl]amino]carbonyl] - (9CI) (CA INDEX NAME)

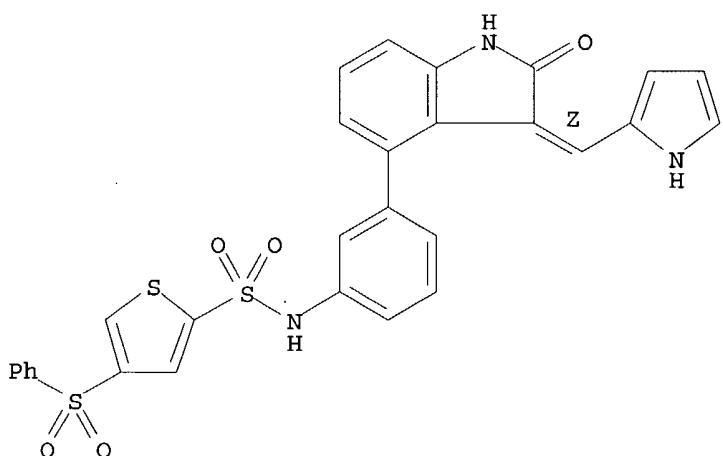
Double bond geometry as shown.



RN 276251-32-2 CAPLUS

CN 2-Thiophenesulfonamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-4-(phenylsulfonyl)- (9CI) (CA INDEX NAME)

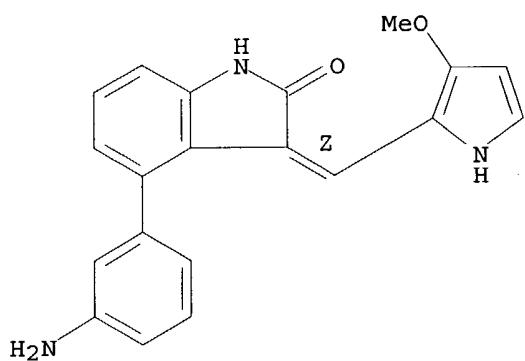
Double bond geometry as shown.



RN 276251-33-3 CAPLUS

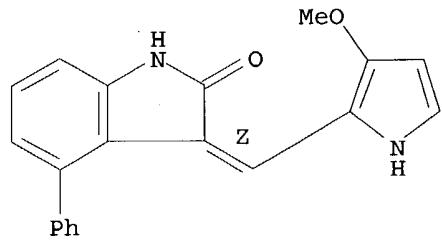
CN 2H-Indol-2-one, 4-(3-aminophenyl)-1,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



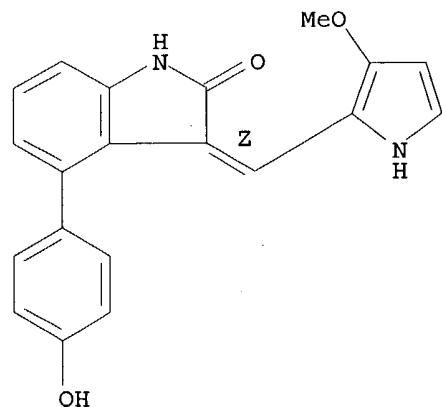
RN 276251-35-5 CAPLUS  
 CN 2H-Indol-2-one, 1,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-4-phenyl-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



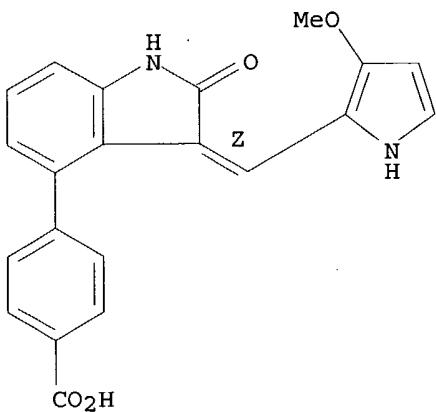
RN 276251-36-6 CAPLUS  
 CN 2H-Indol-2-one, 1,3-dihydro-4-(4-hydroxyphenyl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 276251-37-7 CAPLUS  
 CN Benzoic acid, 4-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-1H-indol-4-yl]- (9CI) (CA INDEX NAME)

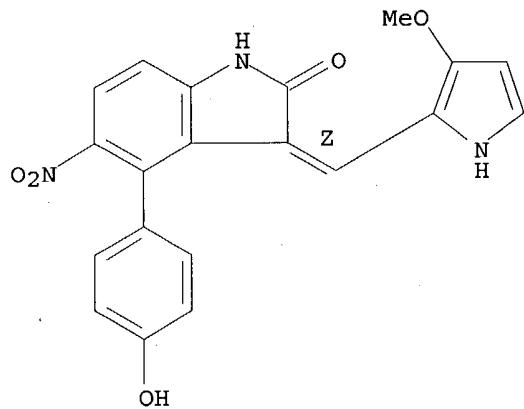
Double bond geometry as shown.



RN 276251-38-8 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(4-hydroxyphenyl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-5-nitro-, (3Z)- (9CI) (CA INDEX NAME)

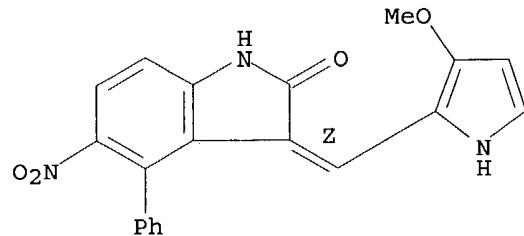
Double bond geometry as shown.



RN 276251-39-9 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-5-nitro-4-phenyl-, (3Z)- (9CI) (CA INDEX NAME)

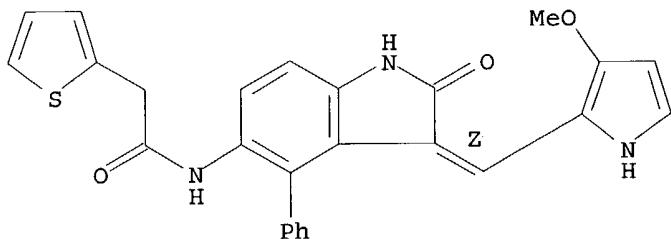
Double bond geometry as shown.



RN 276251-40-2 CAPLUS

CN 2-Thiopheneacetamide, N-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-4-phenyl-1H-indol-5-yl]- (9CI) (CA INDEX NAME)

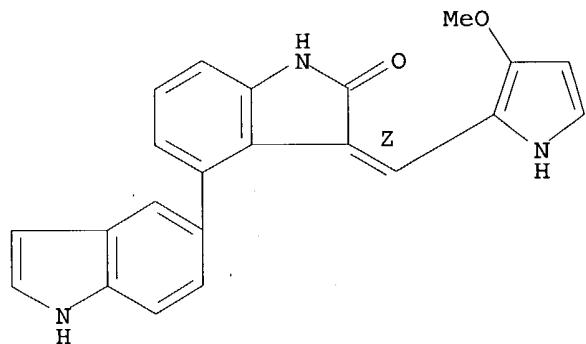
Double bond geometry as shown.



RN 276251-41-3 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(1H-indol-5-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

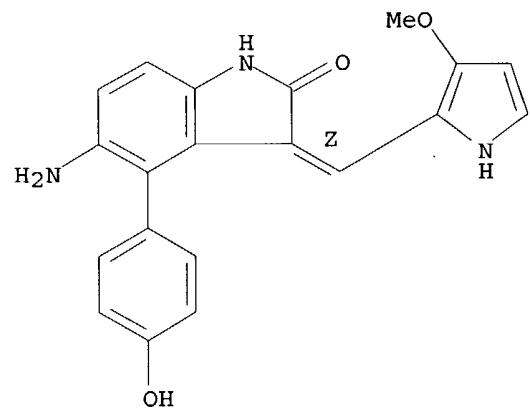
Double bond geometry as shown.



RN 276251-42-4 CAPLUS

CN 2H-Indol-2-one, 5-amino-1,3-dihydro-4-(4-hydroxyphenyl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

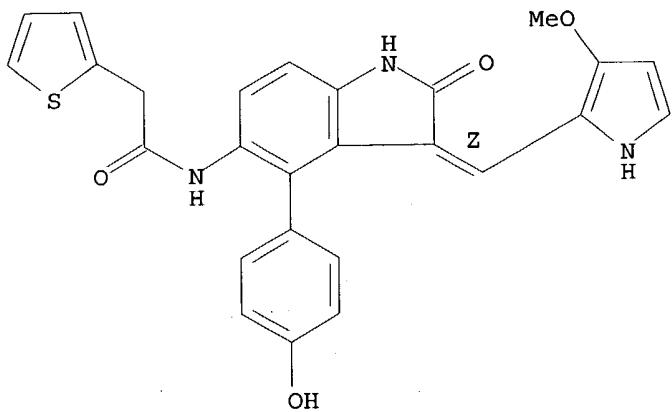
Double bond geometry as shown.



RN 276251-43-5 CAPLUS

CN 2-Thiopheneacetamide, N-[(3Z)-2,3-dihydro-4-(4-hydroxyphenyl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-1H-indol-5-yl]- (9CI) (CA INDEX NAME)

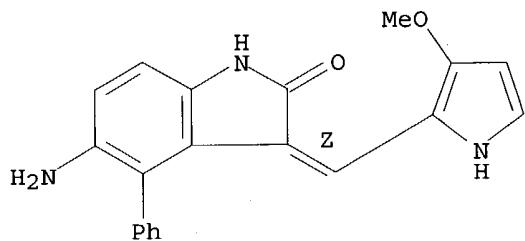
Double bond geometry as shown.



RN 276251-44-6 CAPLUS

CN 2H-Indol-2-one, 5-amino-1,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-4-phenyl-, (3Z)- (9CI) (CA INDEX NAME)

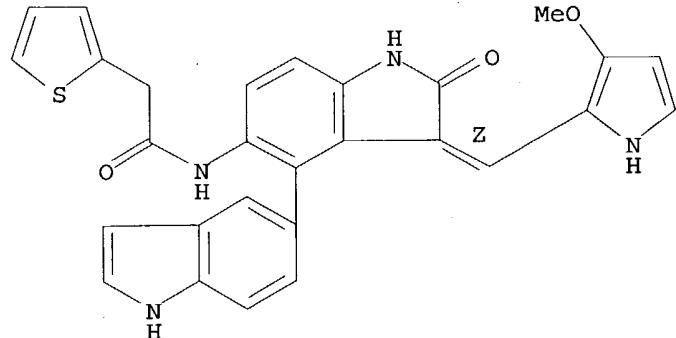
Double bond geometry as shown.



RN 276251-45-7 CAPLUS

CN 2-Thiopheneacetamide, N-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo[4,5'-bi-1H-indol]-5-yl]- (9CI) (CA INDEX NAME)

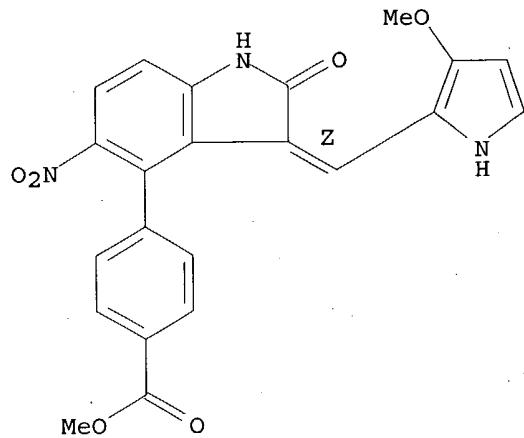
Double bond geometry as shown.



RN 276251-46-8 CAPLUS

CN Benzoic acid, 4-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-5-nitro-2-oxo-1H-indol-4-yl]-, methyl ester (9CI) (CA INDEX NAME)

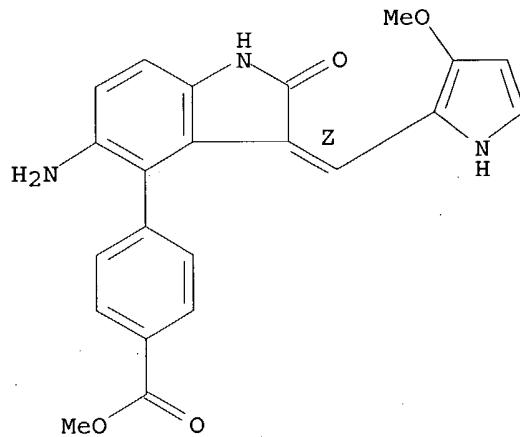
Double bond geometry as shown.



RN 276251-47-9 CAPLUS

CN Benzoic acid, 4-[(3Z)-5-amino-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-1H-indol-4-yl]-, methyl ester (9CI) (CA INDEX NAME)

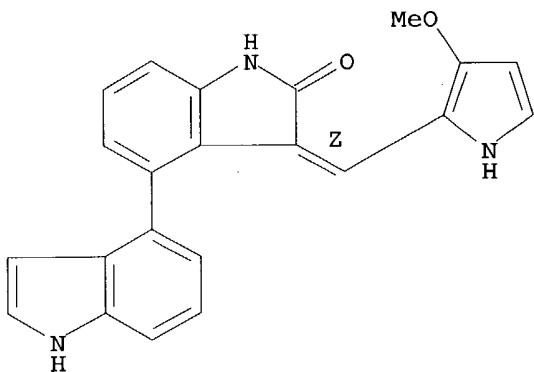
Double bond geometry as shown.



RN 276251-48-0 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(1H-indol-4-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

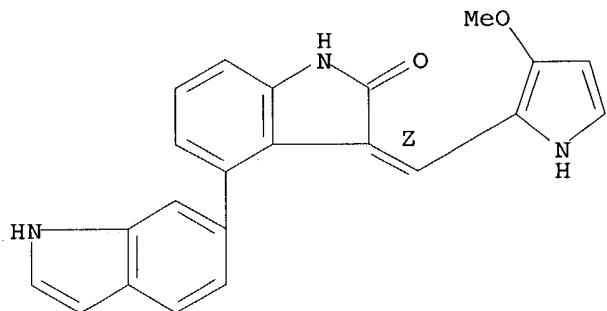
Double bond geometry as shown.



RN 276251-49-1 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(1H-indol-6-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

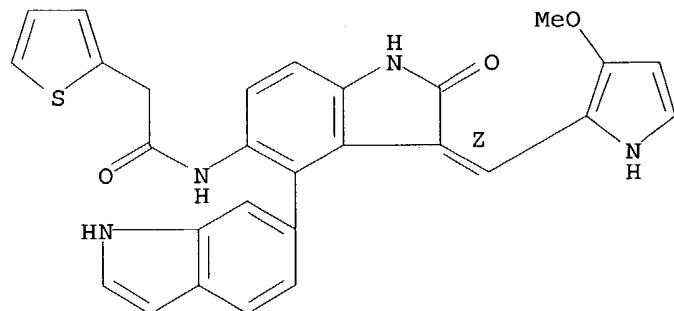
Double bond geometry as shown.



RN 276251-50-4 CAPLUS

CN 2-Thiopheneacetamide, N-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo[4,6'-bi-1H-indol]-5-yl]- (9CI) (CA INDEX NAME)

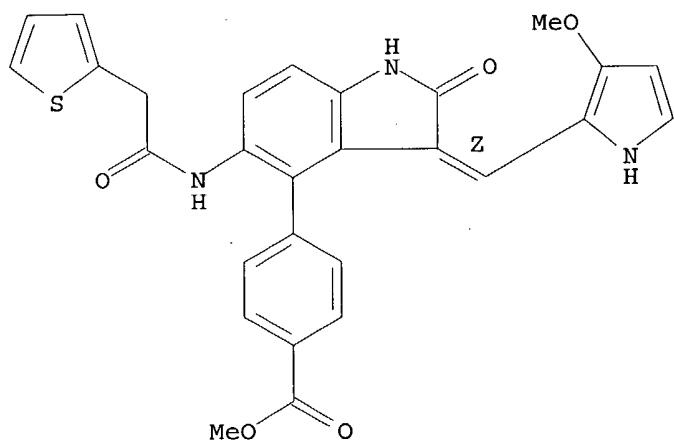
Double bond geometry as shown.



RN 276251-51-5 CAPLUS

CN Benzoic acid, 4-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-5-[(2-thienylacetyl)amino]-1H-indol-4-yl]-, methyl ester (9CI) (CA INDEX NAME)

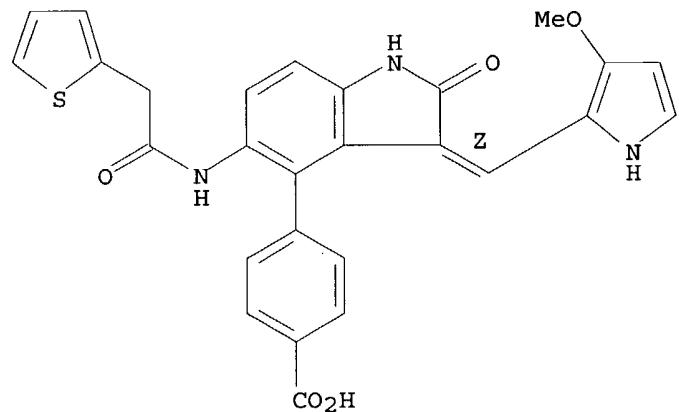
Double bond geometry as shown.



RN 276251-52-6 CAPLUS

CN Benzoic acid, 4-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-5-[(2-thienylacetyl)amino]-1H-indol-4-yl]- (9CI) (CA INDEX NAME)

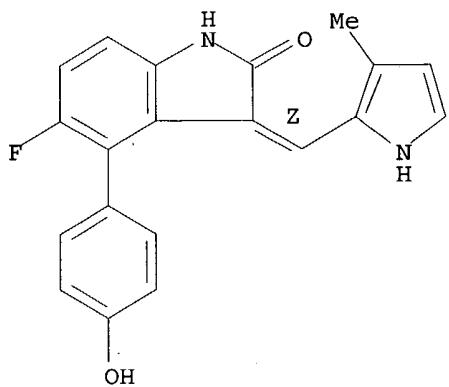
Double bond geometry as shown.



RN 276251-67-3 CAPLUS

CN 2H-Indol-2-one, 5-fluoro-1,3-dihydro-4-(4-hydroxyphenyl)-3-[(3-methyl-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

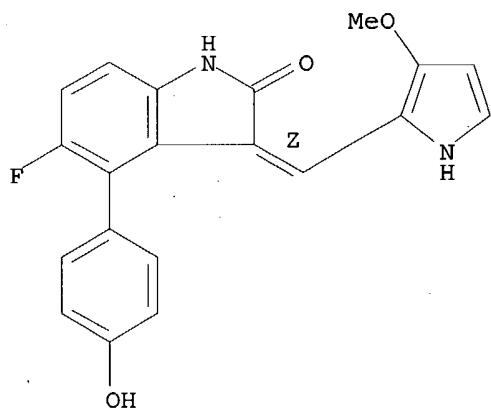
Double bond geometry as shown.



RN 276251-68-4 CAPLUS

CN 2H-Indol-2-one, 5-fluoro-1,3-dihydro-4-(4-hydroxyphenyl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

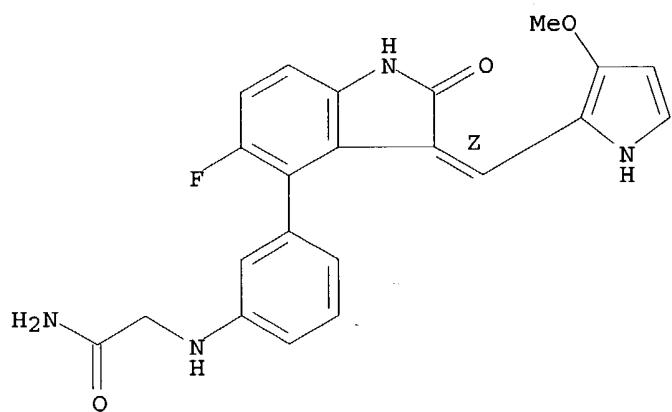
Double bond geometry as shown.



RN 276251-69-5 CAPLUS

CN Acetamide, 2-[[3-[(3Z)-5-fluoro-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-1H-indol-4-yl]phenyl]amino]- (9CI) (CA INDEX NAME)

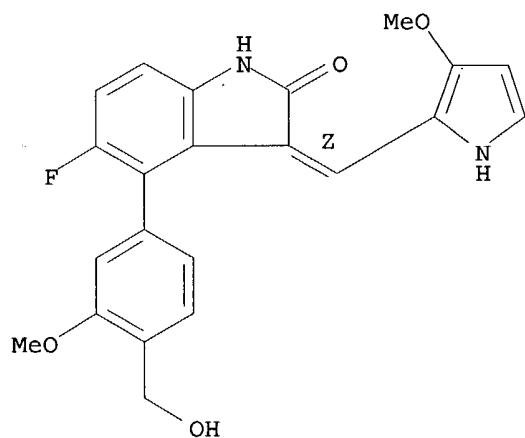
Double bond geometry as shown.



RN 276251-70-8 CAPLUS

CN 2H-Indol-2-one, 5-fluoro-1,3-dihydro-4-[4-(hydroxymethyl)-3-methoxyphenyl]-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

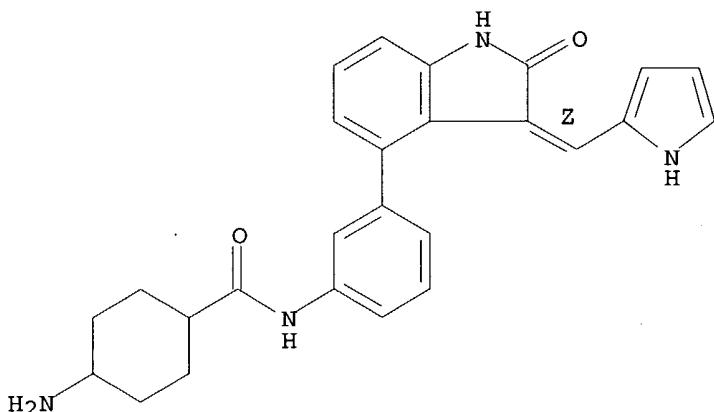
Double bond geometry as shown.



RN 276256-00-9 CAPLUS

CN Cyclohexanecarboxamide, 4-amino-N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



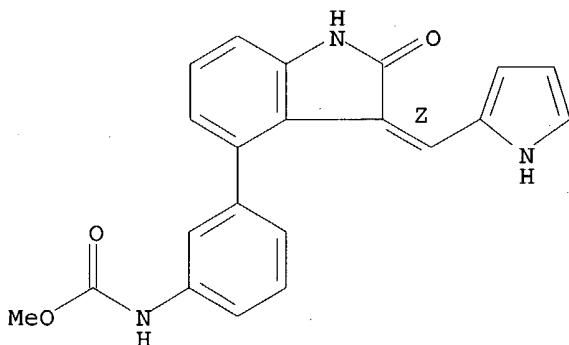
IT 276251-76-4P 276251-79-7P 276251-80-0P  
 276251-81-1P 276251-82-2P 276256-01-0P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)  
 (preparation of 4-aryl-3-(azolylmethylene)-2-oxindoles as inhibitors of JNK protein kinases)

RN 276251-76-4 CAPLUS

CN Carbamic acid, [3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

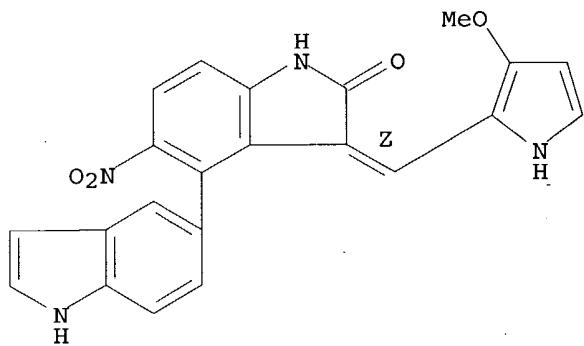
Double bond geometry as shown.



RN 276251-79-7 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(1H-indol-5-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-5-nitro-, (3Z)- (9CI) (CA INDEX NAME)

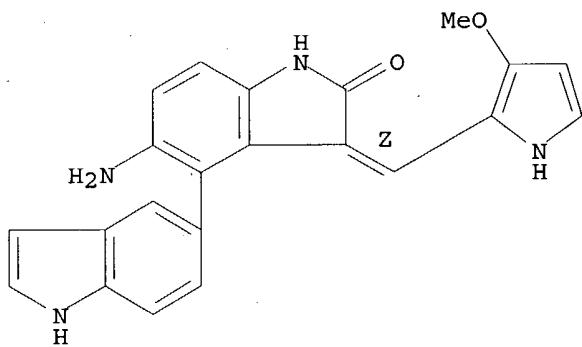
Double bond geometry as shown.



RN 276251-80-0 CAPLUS

CN 2H-Indol-2-one, 5-amino-1,3-dihydro-4-(1H-indol-5-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

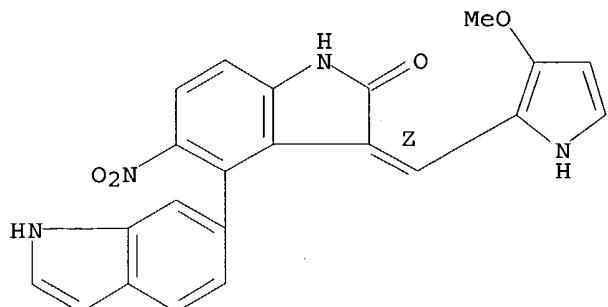
Double bond geometry as shown.



RN 276251-81-1 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(1H-indol-6-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-5-nitro-, (3Z)- (9CI) (CA INDEX NAME)

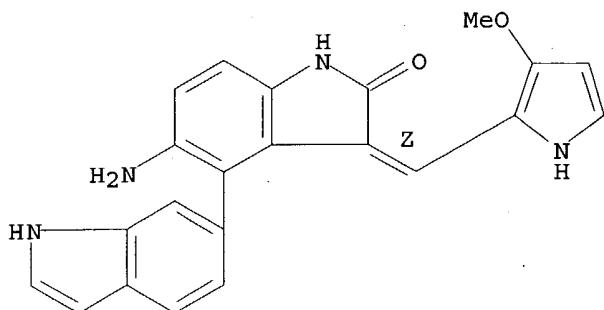
Double bond geometry as shown.



RN 276251-82-2 CAPLUS

CN 2H-Indol-2-one, 5-amino-1,3-dihydro-4-(1H-indol-6-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

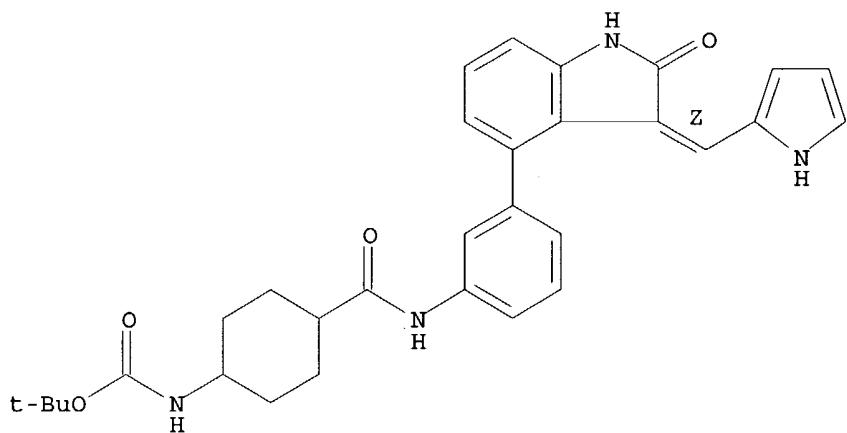
Double bond geometry as shown.



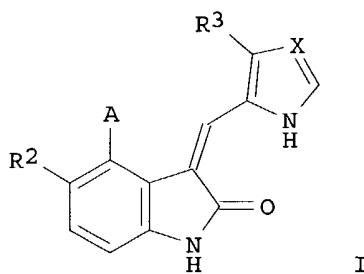
RN 276256-01-0 CAPLUS

CN Carbamic acid, [4-[[[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]amino]carbonyl]cyclohexyl]-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

Double bond geometry as shown.



GI



AB Title compds. [I; A = (substituted) aryl, heteroaryl; R2 = H, halo, OR4, NR6R7, COR4, CO2R4, cyano, NO2, SO2R4, SO2NR6R7, etc.; R3 = H, OR4, COR4, CO2R4, CONR6R7, halo, cyano, NR6R7, perfluoroalkyl, (substituted) alkyl,

etc.; R4 = H, (substituted) alkyl, cycloalkyl, heterocyclyl; R6, R7 = H, (substituted) alkyl, cycloalkyl, COR8, CO2R8, SO2R8, etc.; NR6R7 = (substituted) 3-7 membered ring; R8 = H, (substituted) alkyl, aryl, heteroaryl, cycloalkyl; X = N, CH], were prepared. Thus, (Z)-1,3-dihydro-4-iodo-3-[(1H-pyrrol-2-yl)methylene]-2H-indol-2-one (preparation given) was heated with phenylboronic acid, Pd(OAc)2, Et3N, and tri-O-tolylphosphine in DMF at 100° for 24 h to give 85% (Z)-1,3-dihydro-4-phenyl-3-[(1H-pyrrol-2-yl)methylene]-2H-indol-2-one. Tested I inhibited SAPK with IC50<0.15 μM.

RE.CNT 2 THERE ARE 2 CITED REFERENCES AVAILABLE FOR THIS RECORD  
ALL CITATIONS AVAILABLE IN THE RE FORMAT

DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF,  
 CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

			US 1998-112590P	P	19981217
			US 1999-149028P	P	19990816
BR 9916223	A	20010904	BR 1999-16223		19991209
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EP 1149093	A1	20011031	EP 1999-966933		19991209
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OS MARPAT 133:43433

IT 276250-95-4P 276250-97-6P 276250-98-7P  
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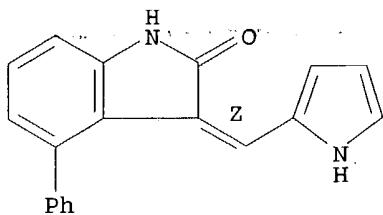
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of 4-aryl-3-(azolylmethylidene)-2-oxindoles as inhibitors of JNK protein kinases)

RN 276250-95-4 CAPLUS

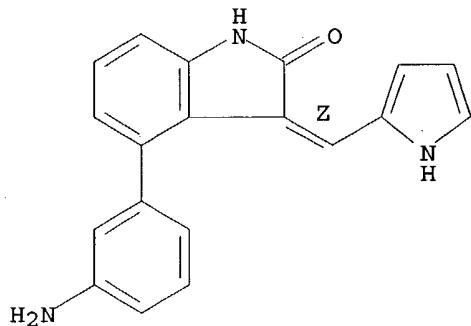
CN 2H-Indol-2-one, 1,3-dihydro-4-phenyl-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



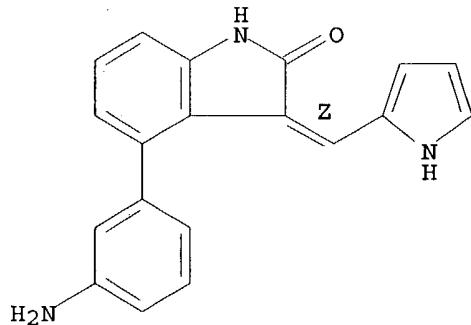
RN 276250-97-6 CAPLUS  
 CN 2H-Indol-2-one, 4-(3-aminophenyl)-1,3-dihydro-3-(1H-pyrrol-2-ylmethylene)-  
 , (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 276250-98-7 CAPLUS  
 CN 2H-Indol-2-one, 4-(3-aminophenyl)-1,3-dihydro-3-(1H-pyrrol-2-ylmethylene)-  
 , monohydrochloride, (3Z)- (9CI) (CA INDEX NAME)

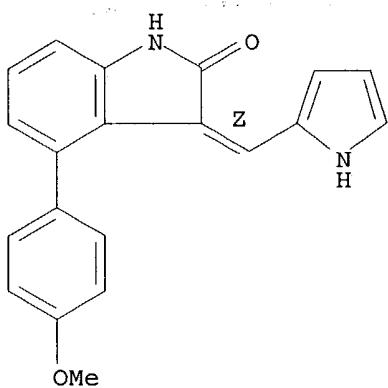
Double bond geometry as shown.



● HCl

RN 276250-99-8 CAPLUS  
 CN 2H-Indol-2-one, 1,3-dihydro-4-(4-methoxyphenyl)-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

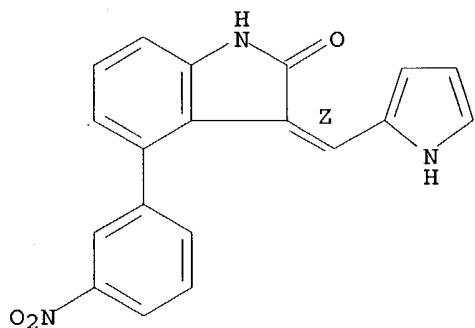
Double bond geometry as shown.



RN 276251-00-4 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(3-nitrophenyl)-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

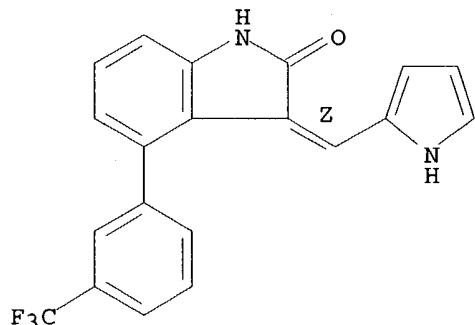
Double bond geometry as shown.



RN 276251-02-6 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-3-(1H-pyrrol-2-ylmethylene)-4-[3-(trifluoromethyl)phenyl]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

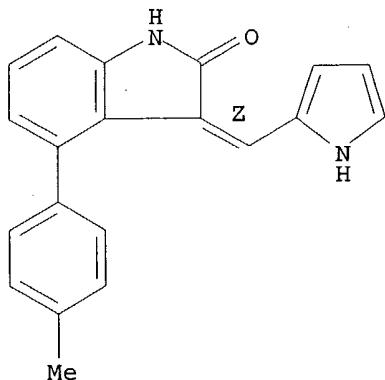


RN 276251-04-8 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(4-methylphenyl)-3-(1H-pyrrol-2-ylmethylene)-

, (3Z)- (9CI) (CA INDEX NAME)

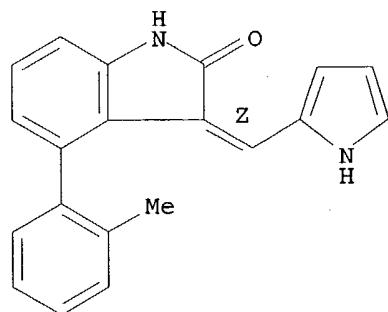
Double bond geometry as shown.



RN 276251-06-0 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(2-methylphenyl)-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

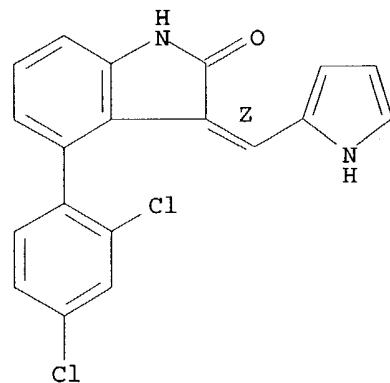
Double bond geometry as shown.



RN 276251-08-2 CAPLUS

CN 2H-Indol-2-one, 4-(2,4-dichlorophenyl)-1,3-dihydro-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

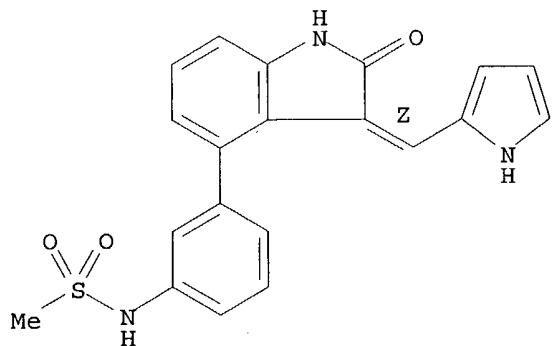
Double bond geometry as shown.



RN 276251-10-6 CAPLUS

CN Methanesulfonamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]- (9CI) (CA INDEX NAME)

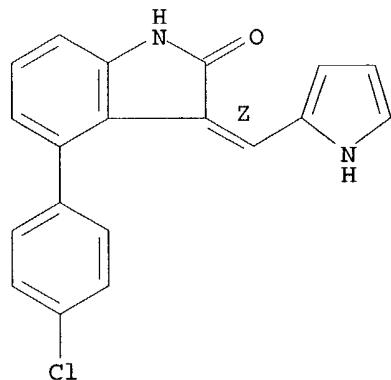
Double bond geometry as shown.



RN 276251-12-8 CAPLUS

CN 2H-Indol-2-one, 4-(4-chlorophenyl)-1,3-dihydro-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

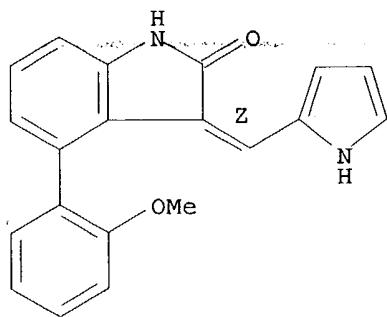
Double bond geometry as shown.



RN 276251-14-0 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(2-methoxyphenyl)-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

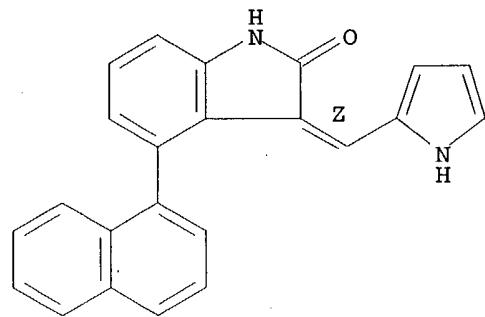
Double bond geometry as shown.



RN 276251-16-2 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(1-naphthalenyl)-3-(1H-pyrrol-2-ylmethylene)-(3Z)-(9CI) (CA INDEX NAME)

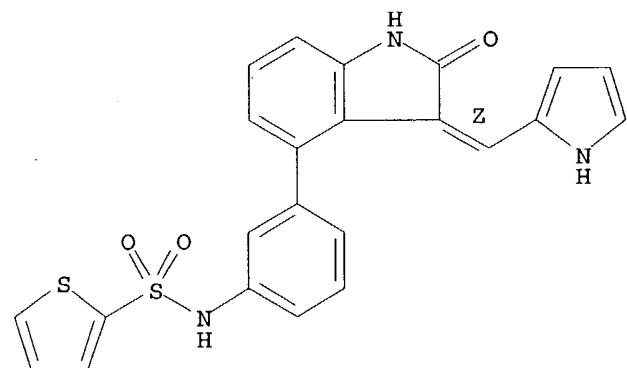
Double bond geometry as shown.



RN 276251-18-4 CAPLUS

CN 2-Thiophenesulfonamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-(9CI) (CA INDEX NAME)

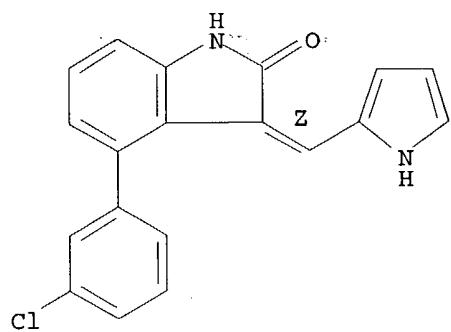
Double bond geometry as shown.



RN 276251-19-5 CAPLUS

CN 2H-Indol-2-one, 4-(3-chlorophenyl)-1,3-dihydro-3-(1H-pyrrol-2-ylmethylene)-(3Z)-(9CI) (CA INDEX NAME)

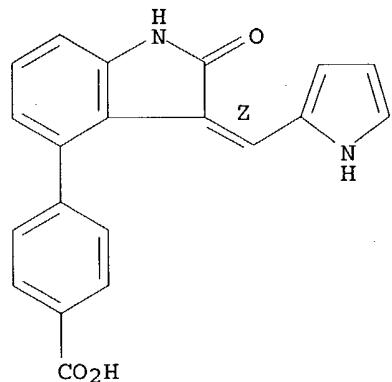
Double bond geometry as shown.



RN 276251-20-8 CAPLUS

CN Benzoic acid, 4-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]- (9CI) (CA INDEX NAME)

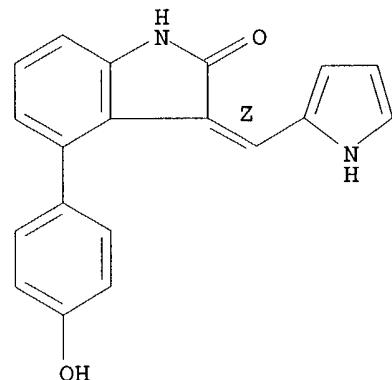
Double bond geometry as shown.



RN 276251-21-9 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(4-hydroxyphenyl)-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

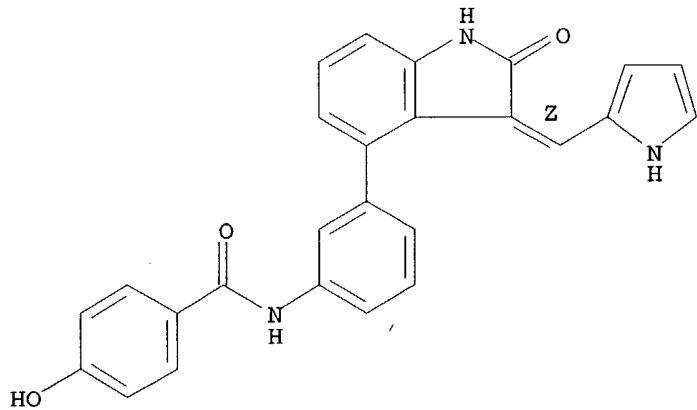
Double bond geometry as shown.



RN 276251-22-0 CAPLUS

CN Benzamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-4-hydroxy- (9CI) (CA INDEX NAME)

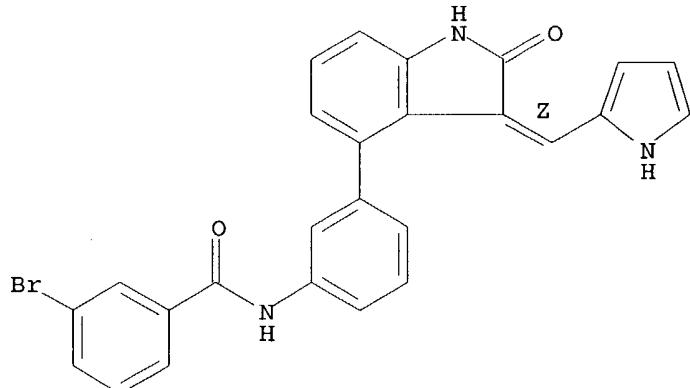
Double bond geometry as shown.



RN 276251-23-1 CAPLUS

CN Benzamide, 3-bromo-N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]- (9CI) (CA INDEX NAME)

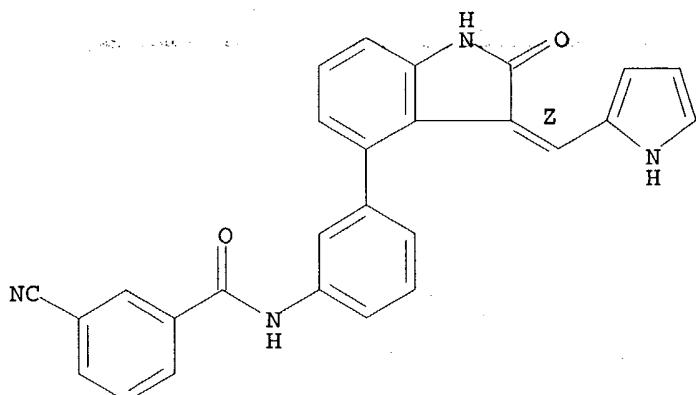
Double bond geometry as shown.



RN 276251-24-2 CAPLUS

CN Benzamide, 3-cyano-N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]- (9CI) (CA INDEX NAME)

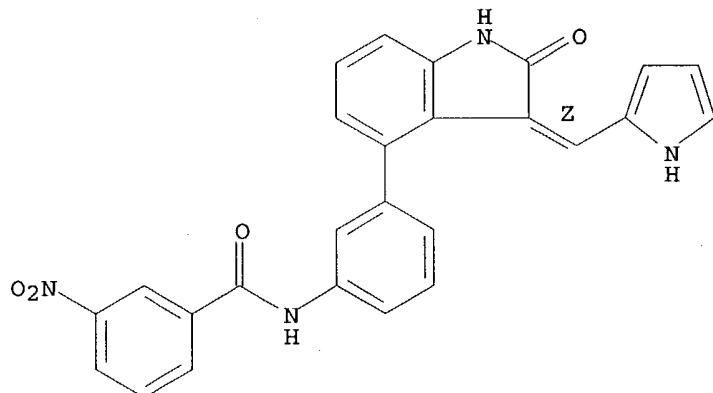
Double bond geometry as shown.



RN 276251-25-3 CAPLUS

CN Benzamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-3-nitro- (9CI) (CA INDEX NAME)

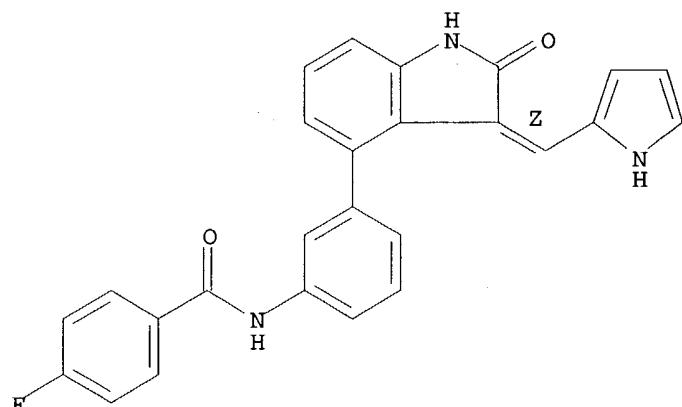
Double bond geometry as shown.



RN 276251-26-4 CAPLUS

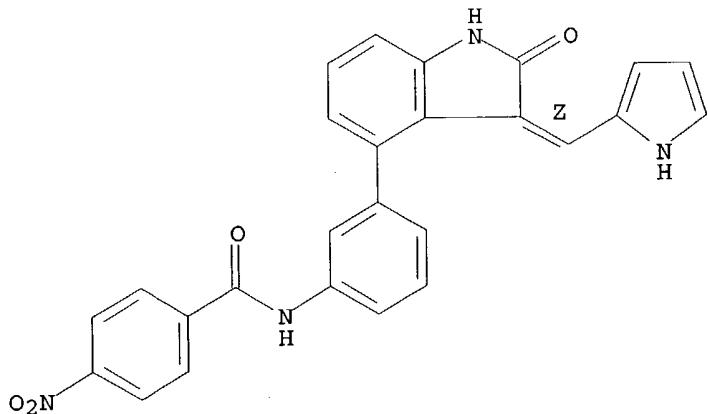
CN Benzamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-4-fluoro- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



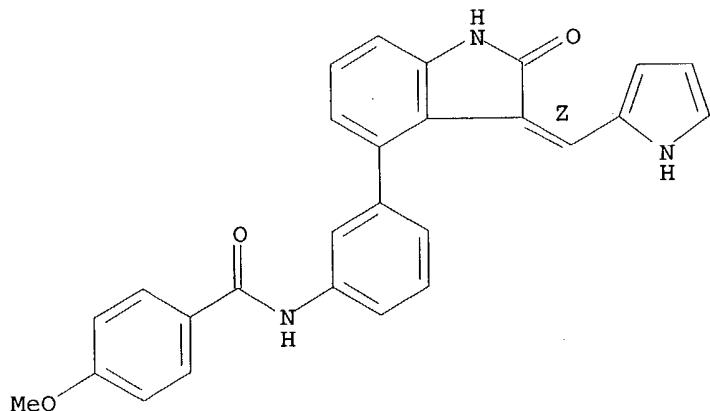
RN 276251-27-5 CAPLUS  
 CN Benzamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-4-nitro- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



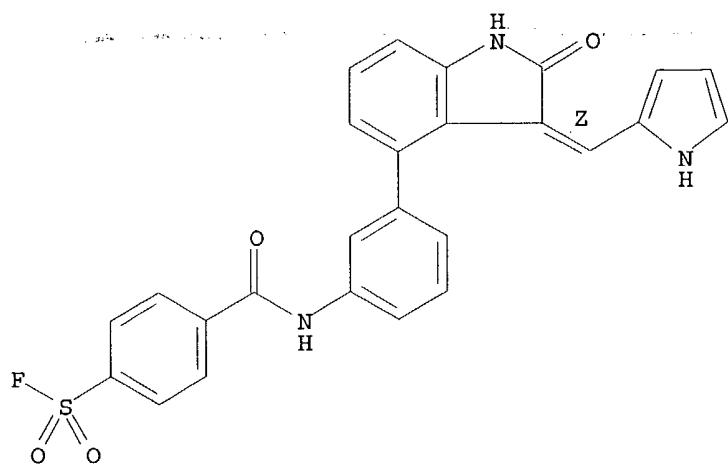
RN 276251-28-6 CAPLUS  
 CN Benzamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-4-methoxy- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 276251-29-7 CAPLUS  
 CN Benzenesulfonyl fluoride, 4-[[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]amino]carbonyl]- (9CI) (CA INDEX NAME)

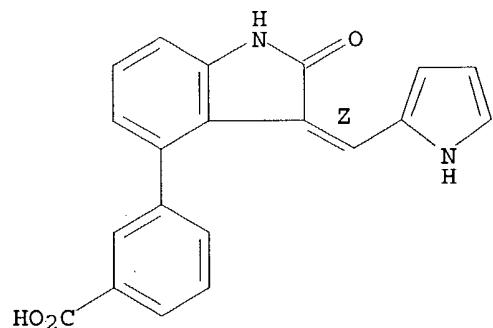
Double bond geometry as shown.



RN 276251-30-0 CAPLUS

CN Benzoic acid, 3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]- (9CI) (CA INDEX NAME)

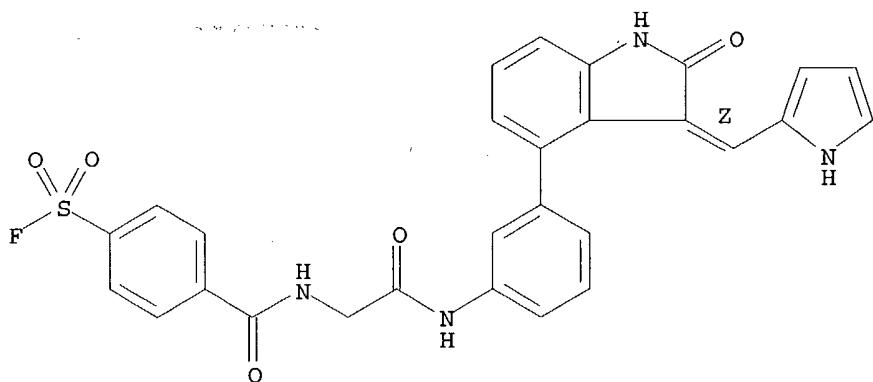
Double bond geometry as shown.



RN 276251-31-1 CAPLUS

CN Benzenesulfonyl fluoride, 4-[[[2-[[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]amino]-2-oxoethyl]amino]carbonyl]- (9CI) (CA INDEX NAME)

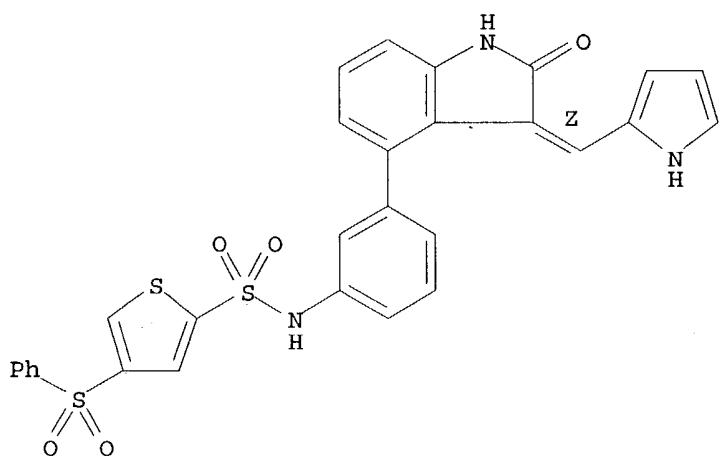
Double bond geometry as shown.



RN 276251-32-2 CAPLUS

CN 2-Thiophenesulfonamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-4-(phenylsulfonyl)- (9CI) (CA INDEX NAME)

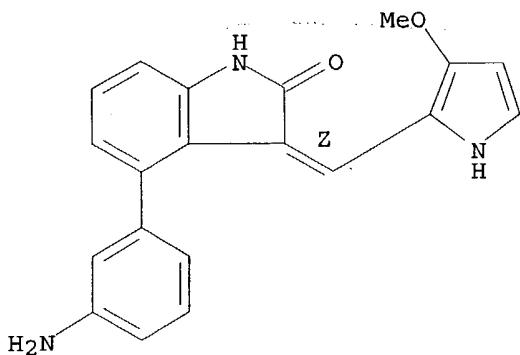
Double bond geometry as shown.



RN 276251-33-3 CAPLUS

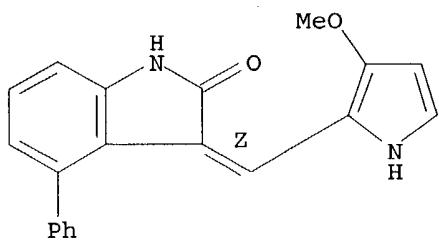
CN 2H-Indol-2-one, 4-(3-aminophenyl)-1,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



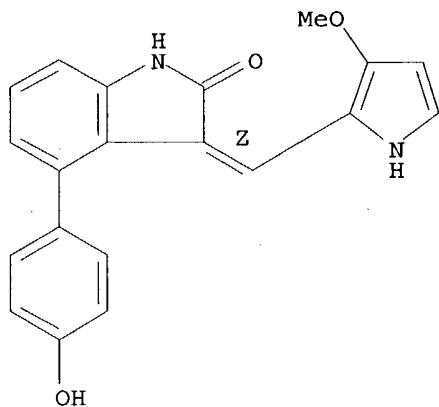
RN 276251-35-5 CAPLUS  
 CN 2H-Indol-2-one, 1,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-4-phenyl-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



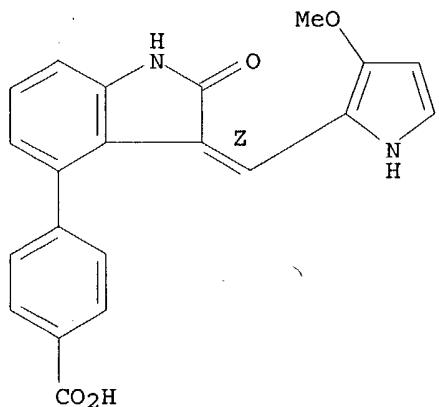
RN 276251-36-6 CAPLUS  
 CN 2H-Indol-2-one, 1,3-dihydro-4-(4-hydroxyphenyl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 276251-37-7 CAPLUS  
 CN Benzoic acid, 4-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-1H-indol-4-yl]- (9CI) (CA INDEX NAME)

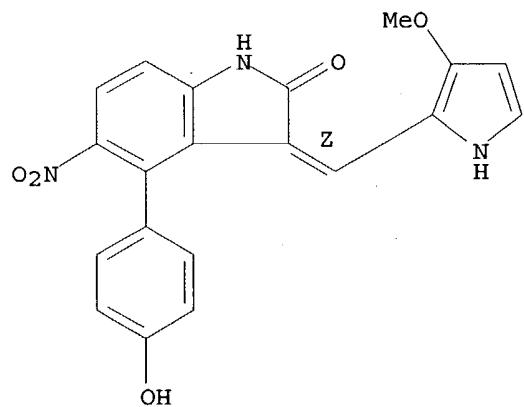
Double bond geometry as shown.



RN 276251-38-8 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(4-hydroxyphenyl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-5-nitro-, (3Z)- (9CI) (CA INDEX NAME)

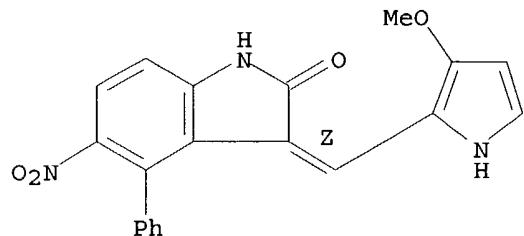
Double bond geometry as shown.



RN 276251-39-9 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-5-nitro-4-phenyl-, (3Z)- (9CI) (CA INDEX NAME)

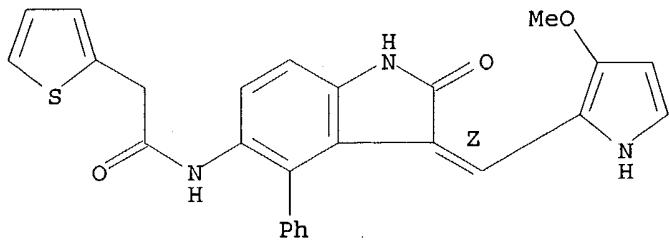
Double bond geometry as shown.



RN 276251-40-2 CAPLUS

CN 2-Thiopheneacetamide, N-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-4-phenyl-1H-indol-5-yl]- (9CI) (CA INDEX NAME)

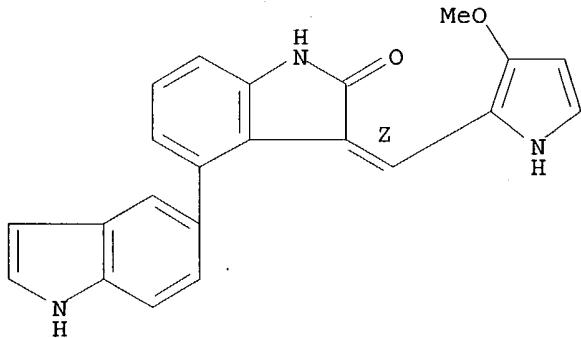
Double bond geometry as shown.



RN 276251-41-3 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(1H-indol-5-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

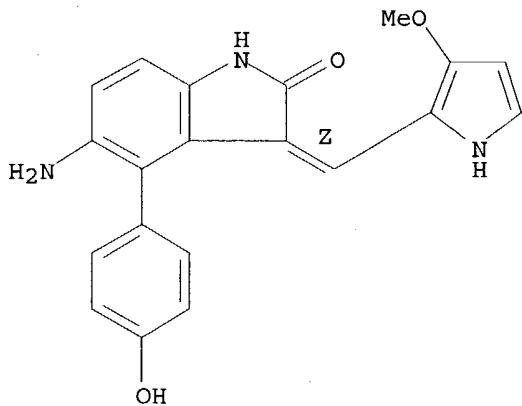
Double bond geometry as shown.



RN 276251-42-4 CAPLUS

CN 2H-Indol-2-one, 5-amino-1,3-dihydro-4-(4-hydroxyphenyl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

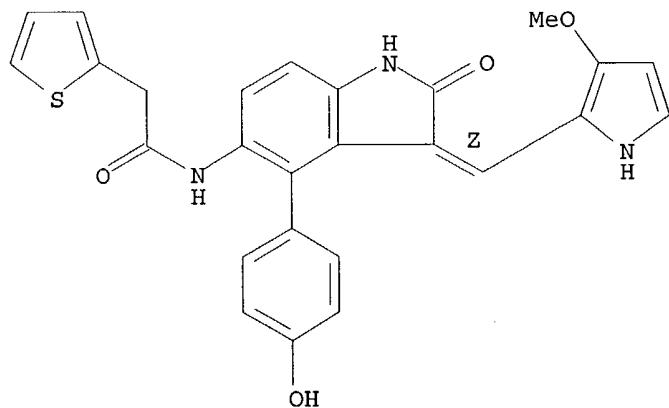
Double bond geometry as shown.



RN 276251-43-5 CAPLUS

2-Thiopheneacetamide, N-[{(3Z)-2,3-dihydro-4-(4-hydroxyphenyl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-1H-indol-5-yl}-(9CI) (CA INDEX NAME)

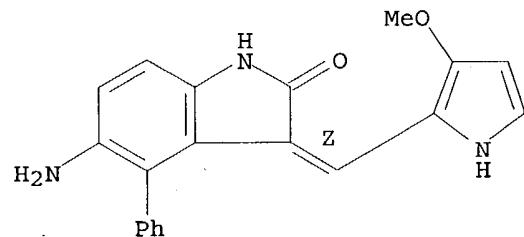
Double bond geometry as shown.



RN 276251-44-6 CAPLUS

CN 2H-Indol-2-one, 5-amino-1,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-4-phenyl-, (3Z)- (9CI) (CA INDEX NAME)

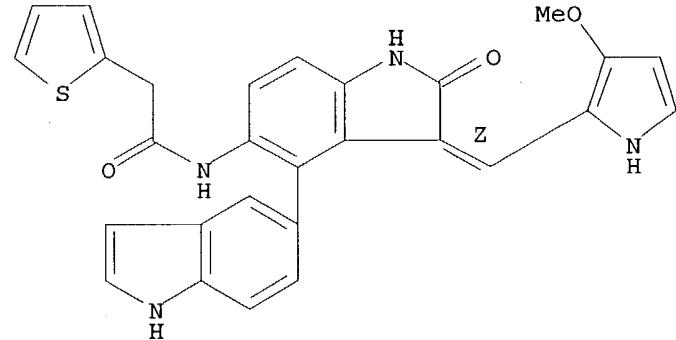
Double bond geometry as shown.



RN 276251-45-7 CAPLUS

CN 2-Thiopheneacetamide, N-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrrol-2-yl)methylene]-2-oxo[4,5'-bi-1H-indol]-5-yl]- (9CI) (CA INDEX NAME)

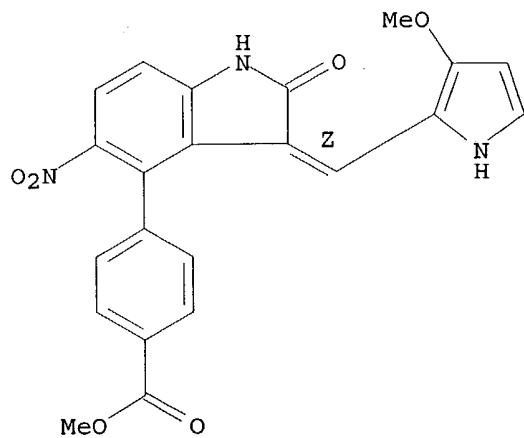
Double bond geometry as shown.



RN 276251-46-8 CAPLUS

CN Benzoic acid, 4-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-5-nitro-2-oxo-1H-indol-4-yl]-, methyl ester (9CI) (CA INDEX NAME)

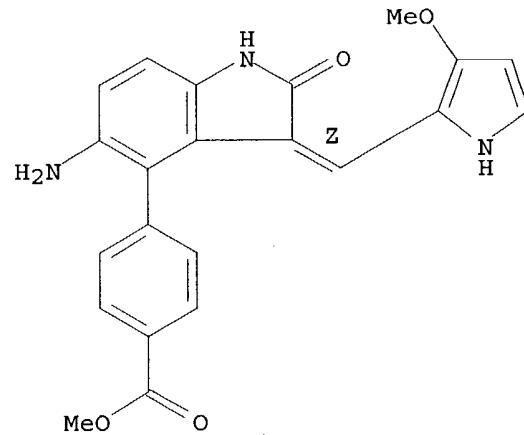
Double bond geometry as shown.



RN 276251-47-9 CAPLUS

CN Benzoic acid, 4-[(3Z)-5-amino-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-1H-indol-4-yl]-, methyl ester (9CI) (CA INDEX NAME)

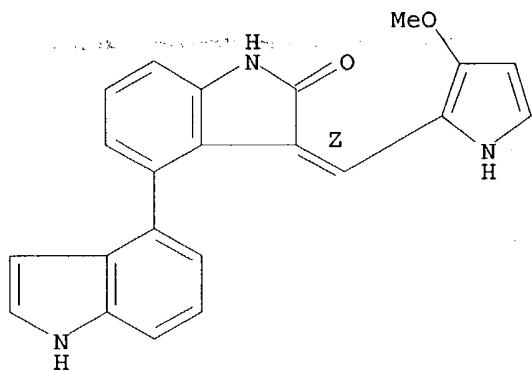
Double bond geometry as shown.



RN 276251-48-0 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(1H-indol-4-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

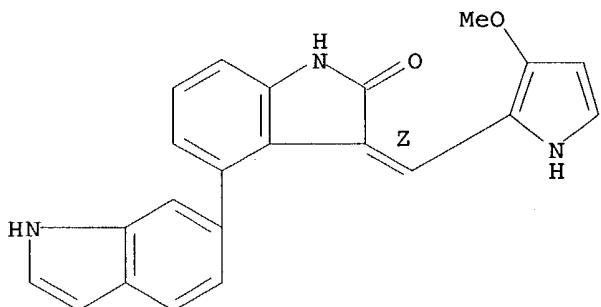
Double bond geometry as shown.



RN 276251-49-1 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(1H-indol-6-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

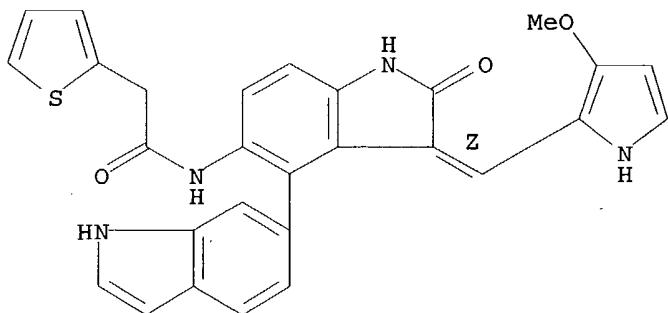
Double bond geometry as shown.



RN 276251-50-4 CAPLUS

CN 2-Thiopheneacetamide, N-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo[4,6'-bi-1H-indol]-5-yl]- (9CI) (CA INDEX NAME)

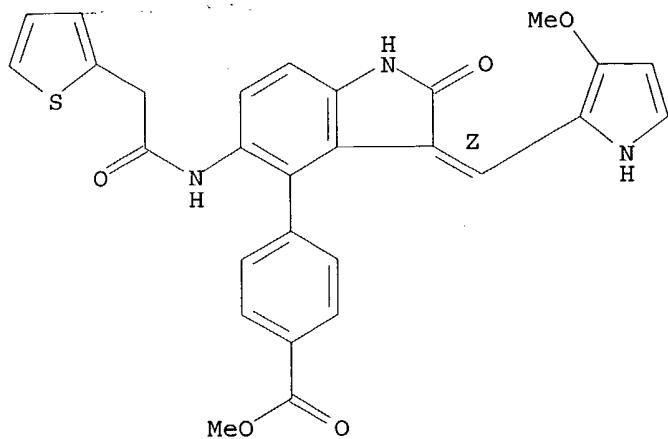
Double bond geometry as shown.



RN 276251-51-5 CAPLUS

CN Benzoic acid, 4-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-5-[(2-thienylacetyl)amino]-1H-indol-4-yl]-, methyl ester (9CI) (CA INDEX NAME)

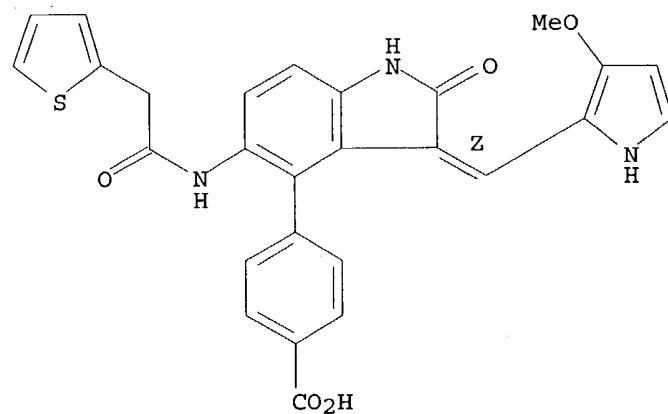
Double bond geometry as shown.



RN 276251-52-6 CAPLUS

CN Benzoic acid, 4-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-5-[(2-thienylacetyl)amino]indol-4-yl]- (9CI) (CA INDEX NAME)

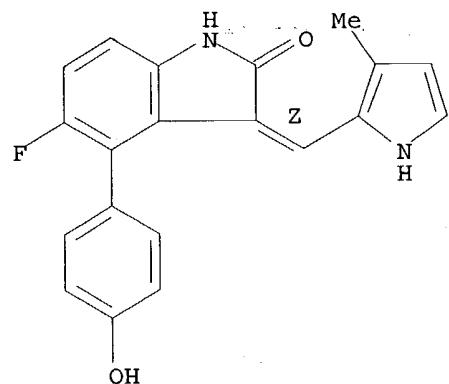
Double bond geometry as shown.



RN 276251-67-3 CAPLUS

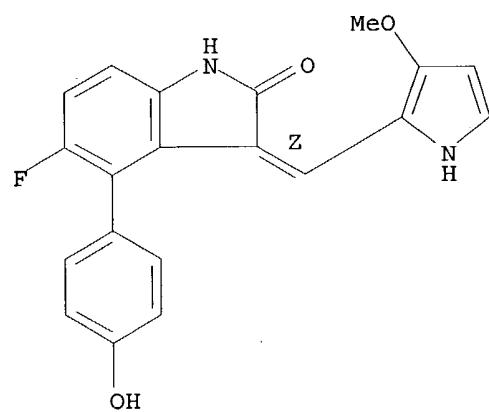
CN 2H-Indol-2-one, 5-fluoro-1,3-dihydro-4-(4-hydroxyphenyl)-3-[(3-methyl-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



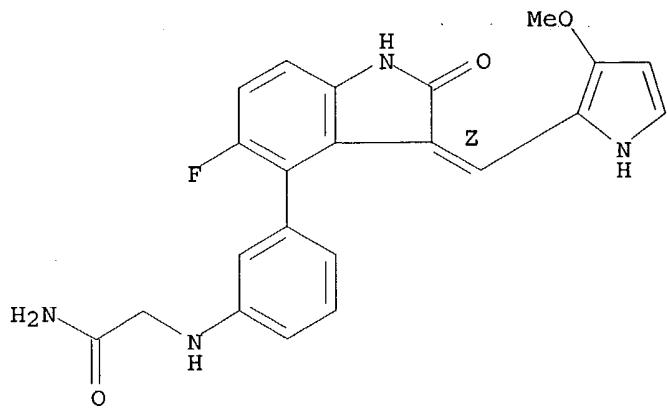
RN 276251-68-4 CAPLUS  
 CN 2H-Indol-2-one, 5-fluoro-1,3-dihydro-4-(4-hydroxyphenyl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 276251-69-5 CAPLUS  
 CN Acetamide, 2-[[3-[(3Z)-5-fluoro-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-1H-indol-4-yl]phenyl]amino]- (9CI) (CA INDEX NAME)

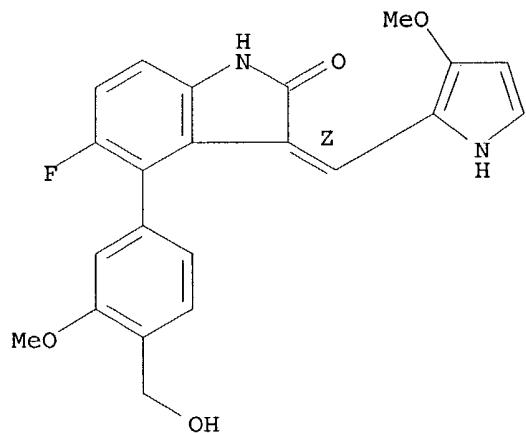
Double bond geometry as shown.



RN 276251-70-8 CAPLUS

CN 2H-Indol-2-one, 5-fluoro-1,3-dihydro-4-[4-(hydroxymethyl)-3-methoxyphenyl]-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

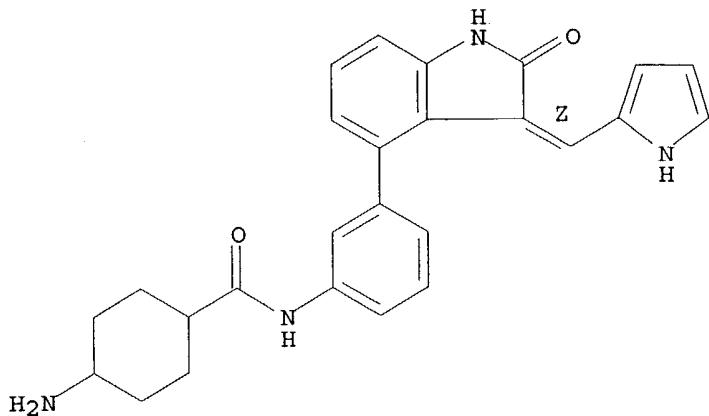
Double bond geometry as shown.



RN 276256-00-9 CAPLUS

CN Cyclohexanecarboxamide, 4-amino-N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



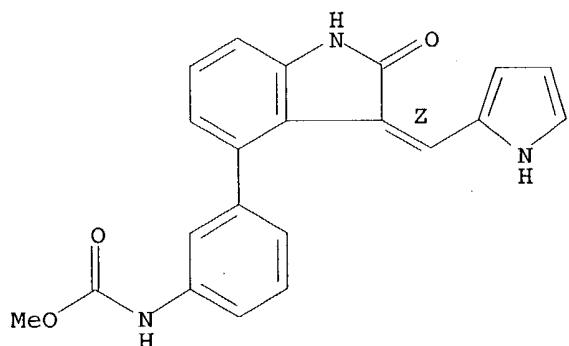
IT 276251-76-4P 276251-79-7P 276251-80-0P  
276251-81-1P 276251-82-2P 276256-01-0P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)  
(preparation of 4-aryl-3-(azolylmethylidene)-2-oxindoles as inhibitors of JNK protein kinases)

RN 276251-76-4 CAPLUS

CN Carbamic acid, [3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

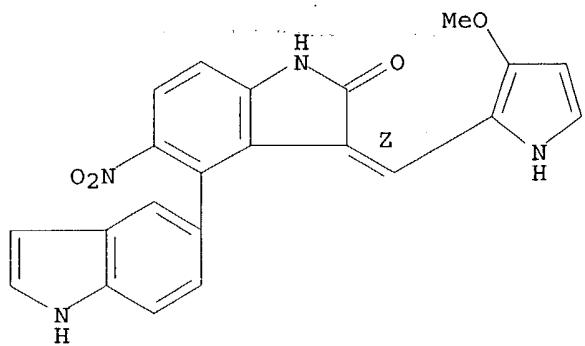
Double bond geometry as shown.



RN 276251-79-7 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(1H-indol-5-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-5-nitro-, (3Z)- (9CI) (CA INDEX NAME)

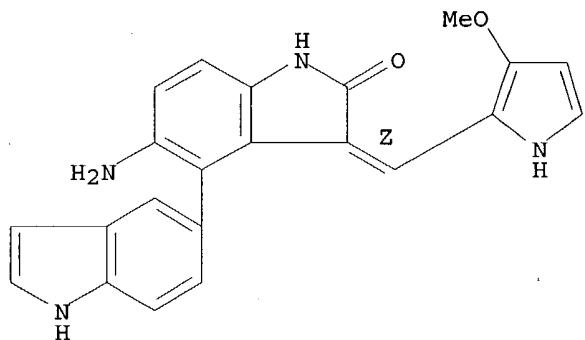
Double bond geometry as shown.



RN 276251-80-0 CAPLUS

CN 2H-Indol-2-one, 5-amino-1,3-dihydro-4-(1H-indol-5-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

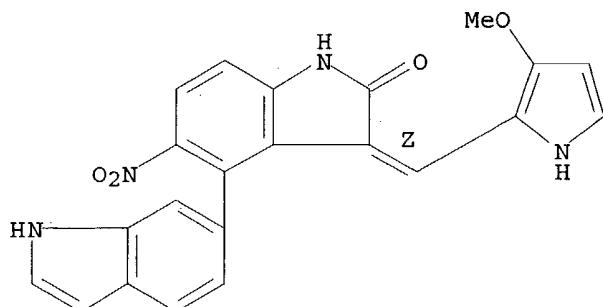
Double bond geometry as shown.



RN 276251-81-1 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(1H-indol-6-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-5-nitro-, (3Z)- (9CI) (CA INDEX NAME)

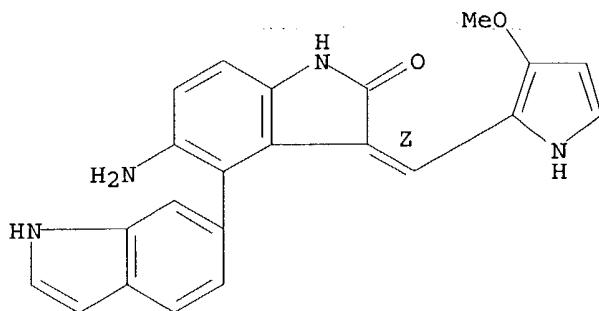
Double bond geometry as shown.



RN 276251-82-2 CAPLUS

CN 2H-Indol-2-one, 5-amino-1,3-dihydro-4-(1H-indol-6-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

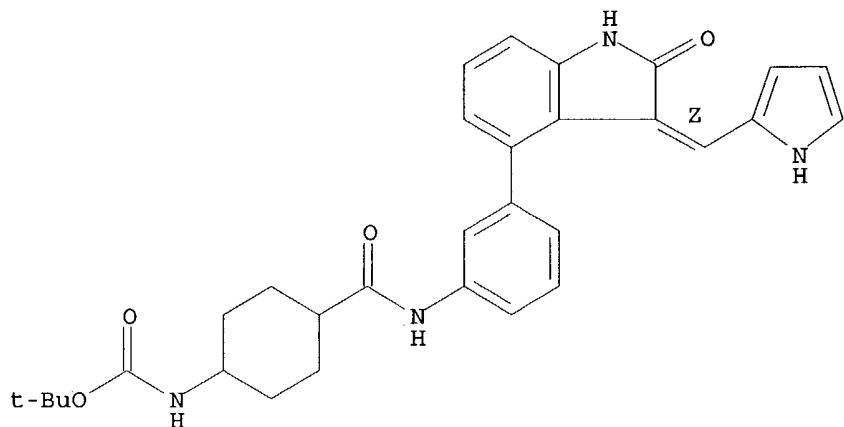
Double bond geometry as shown.



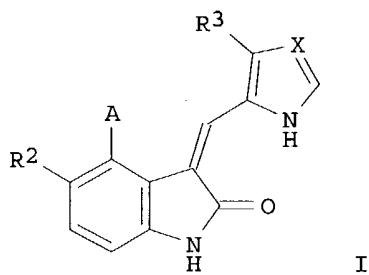
RN 276256-01-0 CAPLUS

CN Carbamic acid, [4-[[[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]amino]carbonyl]cyclohexyl]-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

Double bond geometry as shown.



GI



AB Title compds. [I; A = (substituted) aryl, heteroaryl; R2 = H, halo, OR4, NR6R7, COR4, CO2R4, cyano, NO2, SO2R4, SO2NR6R7, etc.; R3 = H, OR4, COR4, CO2R4, CONR6R7, halo, cyano, NR6R7, perfluoroalkyl, (substituted) alkyl,

etc.; R4 = H, (substituted) alkyl, cycloalkyl, heterocyclyl; R6, R7 = H, (substituted) alkyl, cycloalkyl, COR8, CO2R8, SO2R8, etc.; NR6R7 = (substituted) 3-7 membered ring; R8 = H, (substituted) alkyl, aryl, heteroaryl, cycloalkyl; X = N, CH], were prepared. Thus, (Z)-1,3-dihydro-4-iodo-3-[(1H-pyrrol-2-yl)methylene]-2H-indol-2-one (preparation given) was heated with phenylboronic acid, Pd(OAc)<sub>2</sub>, Et<sub>3</sub>N, and tri-O-tolylphosphine in DMF at 100° for 24 h to give 85% (Z)-1,3-dihydro-4-phenyl-3-[(1H-pyrrol-2-yl)methylene]-2H-indol-2-one. Tested I inhibited SAPK with IC50<0.15 μM.

RE.CNT 2 THERE ARE 2 CITED REFERENCES AVAILABLE FOR THIS RECORD  
ALL CITATIONS AVAILABLE IN THE RE FORMAT

DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF,  
CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

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			US 1999-149028P	P	19990816
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			WO 1999-EP9673	W	19991209
EP 1149093	A1	20011031	EP 1999-966933		19991209
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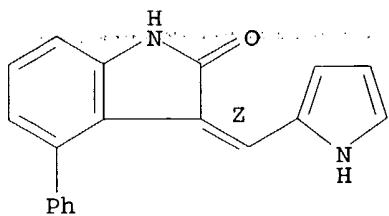
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IT 276250-95-4P 276250-97-6P 276250-98-7P  
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276251-23-1P 276251-24-2P 276251-25-3P  
276251-26-4P 276251-27-5P 276251-28-6P  
276251-29-7P 276251-30-0P 276251-31-1P  
276251-32-2P 276251-33-3P 276251-35-5P  
276251-36-6P 276251-37-7P 276251-38-8P  
276251-39-9P 276251-40-2P 276251-41-3P  
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276251-45-7P 276251-46-8P 276251-47-9P  
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276251-68-4P 276251-69-5P 276251-70-8P  
276256-00-9P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)  
(preparation of 4-aryl-3-(azolylmethylidene)-2-oxindoles as inhibitors of JNK protein kinases)

RN 276250-95-4 CAPLUS  
CN 2H-Indol-2-one, 1,3-dihydro-4-phenyl-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

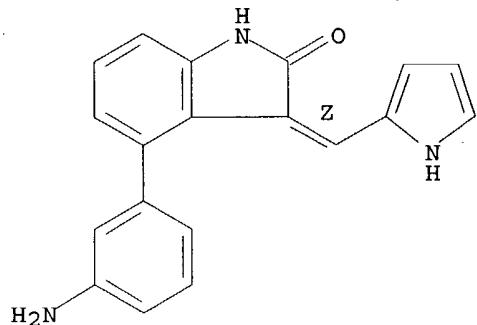
Double bond geometry as shown.



RN 276250-97-6 CAPLUS

CN 2H-Indol-2-one, 4-(3-aminophenyl)-1,3-dihydro-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

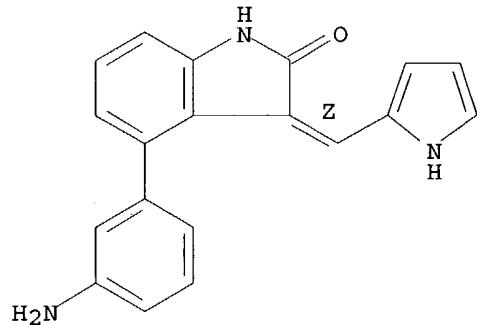
Double bond geometry as shown.



RN 276250-98-7 CAPLUS

CN 2H-Indol-2-one, 4-(3-aminophenyl)-1,3-dihydro-3-(1H-pyrrol-2-ylmethylene)-, monohydrochloride, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

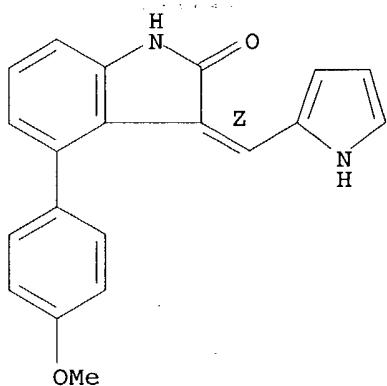


● HCl

RN 276250-99-8 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(4-methoxyphenyl)-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

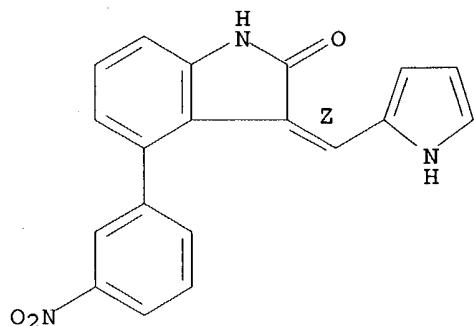
Double bond geometry as shown.



RN 276251-00-4 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(3-nitrophenyl)-3-(1H-pyrrol-2-ylmethylene)-  
(3Z)- (9CI) (CA INDEX NAME)

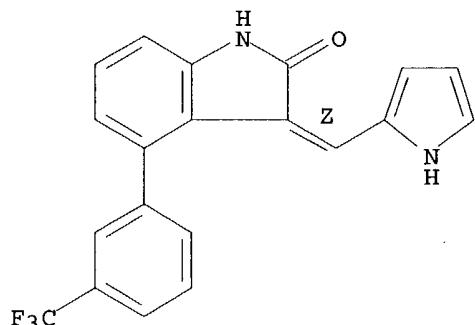
Double bond geometry as shown.



RN 276251-02-6 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-3-(1H-pyrrol-2-ylmethylene)-4-[3-(trifluoromethyl)phenyl]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

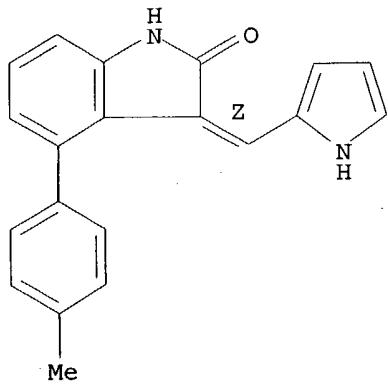


RN 276251-04-8 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(4-methylphenyl)-3-(1H-pyrrol-2-ylmethylene)-

, (3Z)- (9CI) (CA INDEX NAME)

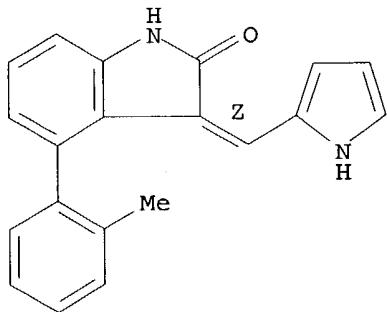
Double bond geometry as shown.



RN 276251-06-0 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(2-methylphenyl)-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

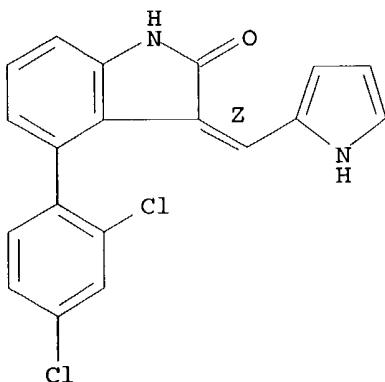
Double bond geometry as shown.



RN 276251-08-2 CAPLUS

CN 2H-Indol-2-one, 4-(2,4-dichlorophenyl)-1,3-dihydro-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

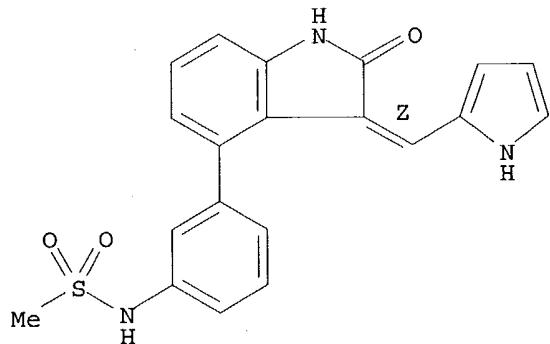
Double bond geometry as shown.



RN 276251-10-6 CAPLUS

CN Methanesulfonamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]- (9CI) (CA INDEX NAME)

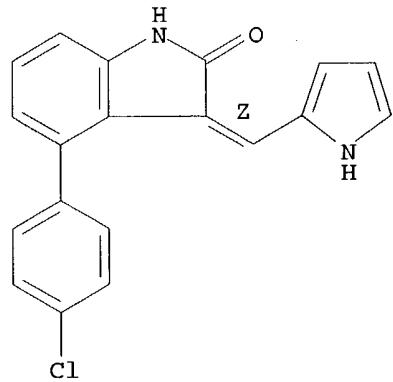
Double bond geometry as shown.



RN 276251-12-8 CAPLUS

CN 2H-Indol-2-one, 4-(4-chlorophenyl)-1,3-dihydro-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

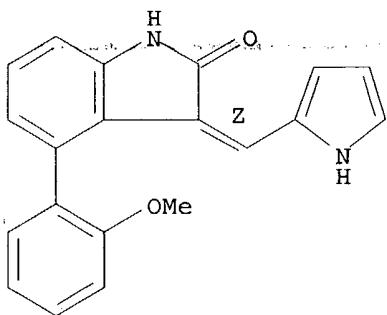
Double bond geometry as shown.



RN 276251-14-0 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(2-methoxyphenyl)-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

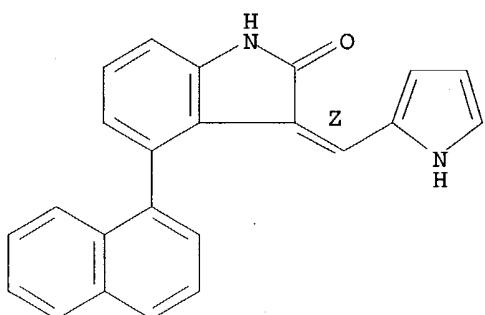
Double bond geometry as shown.



RN 276251-16-2 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(1-naphthalenyl)-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

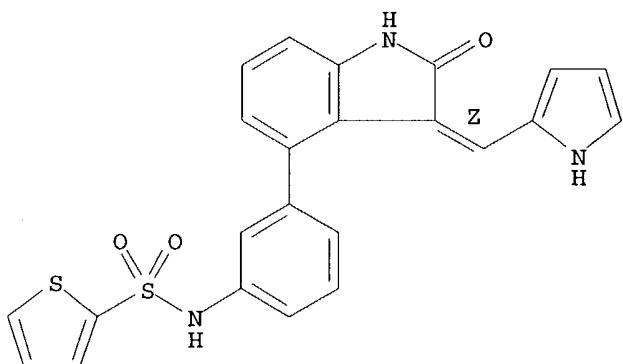
Double bond geometry as shown.



RN 276251-18-4 CAPLUS

CN 2-Thiophenesulfonamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]- (9CI) (CA INDEX NAME)

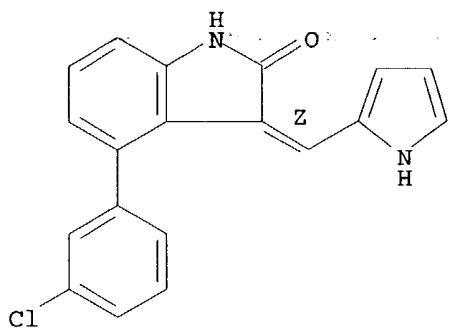
Double bond geometry as shown.



RN 276251-19-5 CAPLUS

CN 2H-Indol-2-one, 4-(3-chlorophenyl)-1,3-dihydro-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

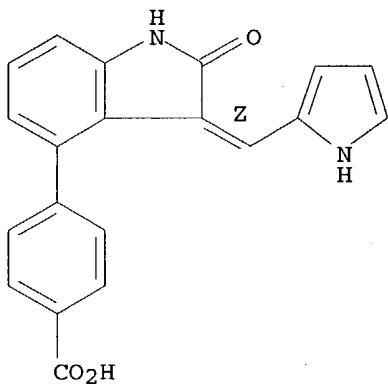
Double bond geometry as shown.



RN 276251-20-8 CAPLUS

CN Benzoic acid, 4-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]- (9CI) (CA INDEX NAME)

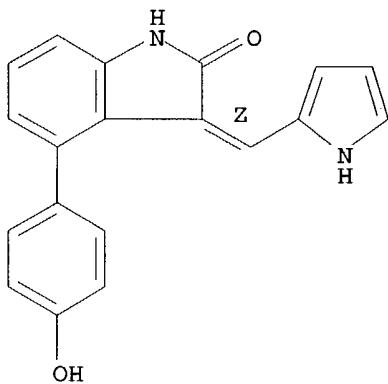
Double bond geometry as shown.



RN 276251-21-9 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(4-hydroxyphenyl)-3-(1H-pyrrol-2-ylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

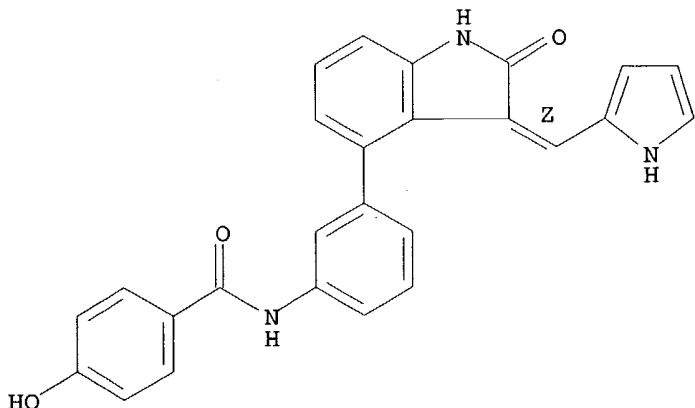
Double bond geometry as shown.



RN 276251-22-0 CAPLUS

CN Benzamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-4-hydroxy- (9CI) (CA INDEX NAME)

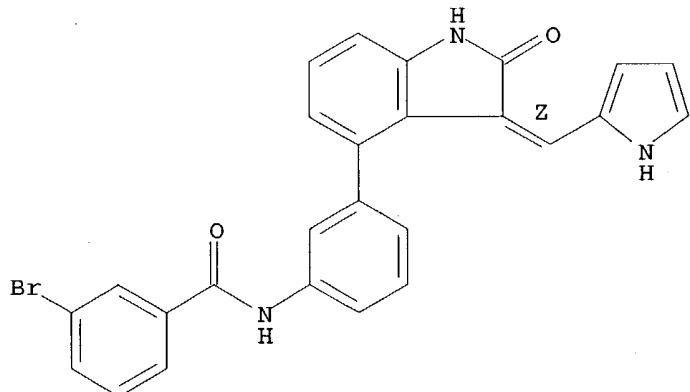
Double bond geometry as shown.



RN 276251-23-1 CAPLUS

CN Benzamide, 3-bromo-N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]- (9CI) (CA INDEX NAME)

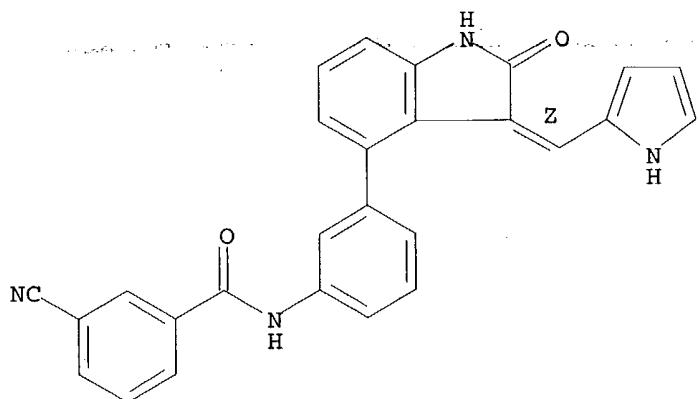
Double bond geometry as shown.



RN 276251-24-2 CAPLUS

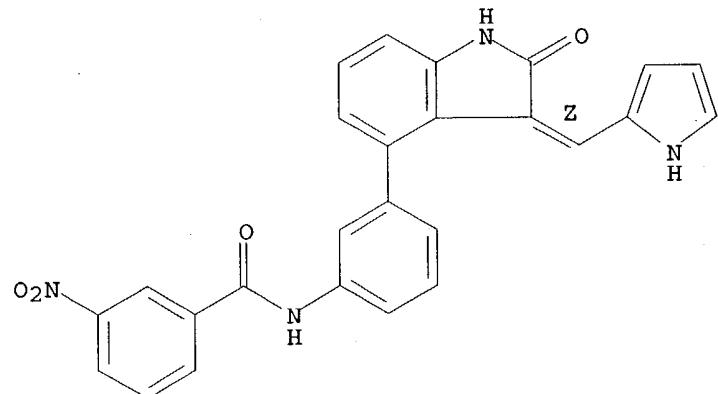
CN Benzamide, 3-cyano-N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



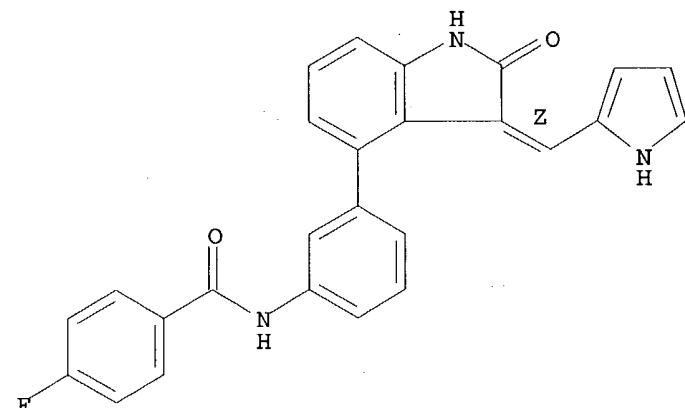
RN 276251-25-3 CAPLUS  
 CN Benzamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-3-nitro- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 276251-26-4 CAPLUS  
 CN Benzamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-4-fluoro- (9CI) (CA INDEX NAME)

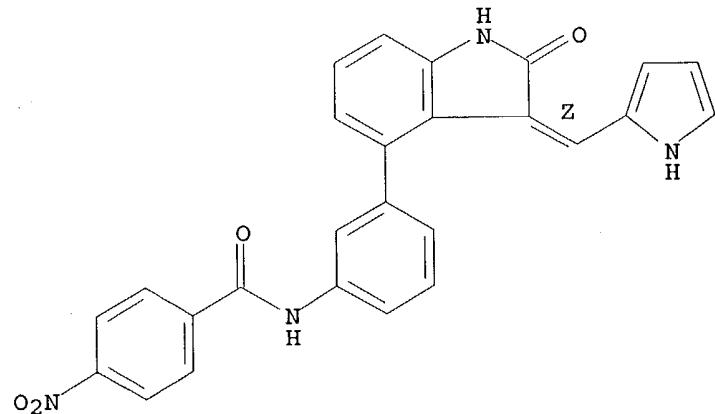
Double bond geometry as shown.



RN 276251-27-5 CAPLUS

CN Benzamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-4-nitro- (9CI) (CA INDEX NAME)

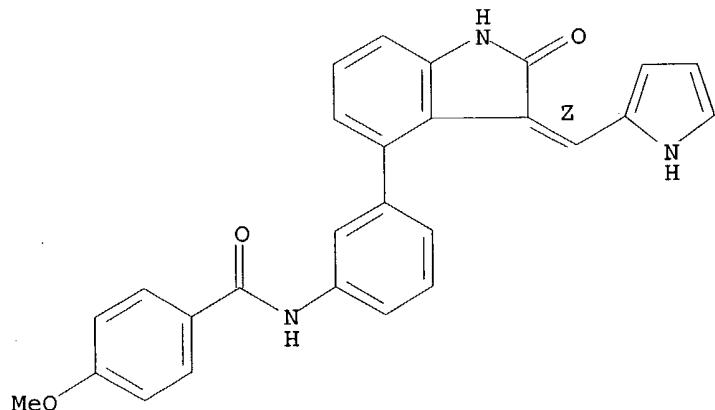
Double bond geometry as shown.



RN 276251-28-6 CAPLUS

CN Benzamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-4-methoxy- (9CI) (CA INDEX NAME)

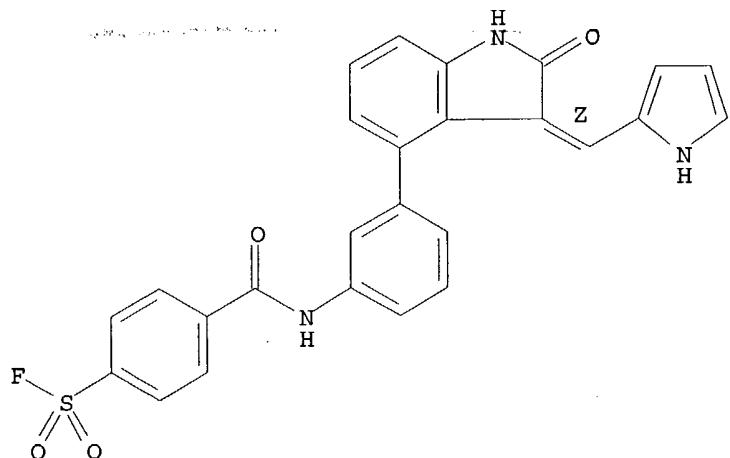
Double bond geometry as shown.



RN 276251-29-7 CAPLUS

CN Benzenesulfonyl fluoride, 4-[[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]amino]carbonyl- (9CI) (CA INDEX NAME)

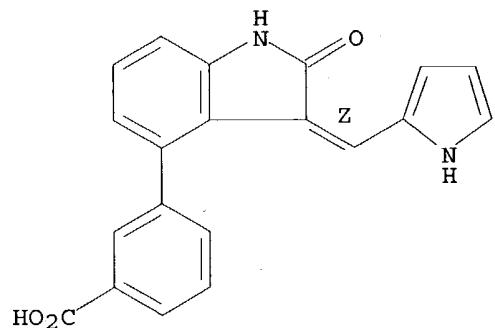
Double bond geometry as shown.



RN 276251-30-0 CAPLUS

CN Benzoic acid, 3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]- (9CI) (CA INDEX NAME)

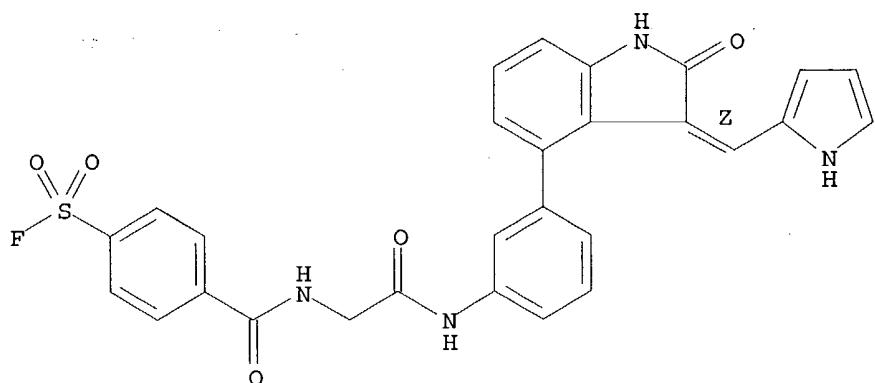
Double bond geometry as shown.



RN 276251-31-1 CAPLUS

CN Benzenesulfonyl fluoride, 4-[[[2-[[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]amino]-2-oxoethyl]amino]carbonyl]- (9CI) (CA INDEX NAME)

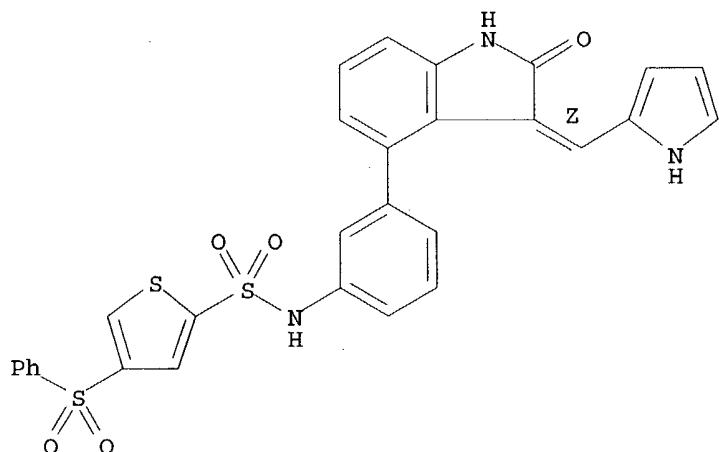
Double bond geometry as shown.



RN 276251-32-2 CAPLUS

CN 2-Thiophenesulfonamide, N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-4-(phenylsulfonyl)- (9CI) (CA INDEX NAME)

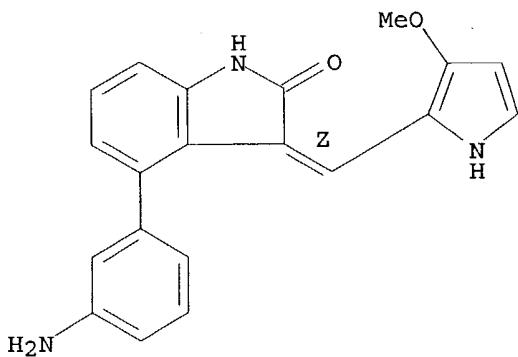
Double bond geometry as shown.



RN 276251-33-3 CAPLUS

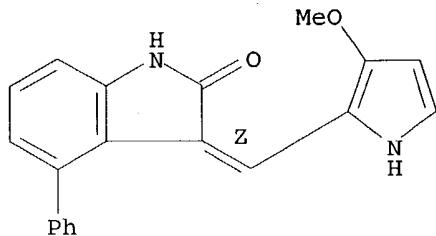
CN 2H-Indol-2-one, 4-(3-aminophenyl)-1,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



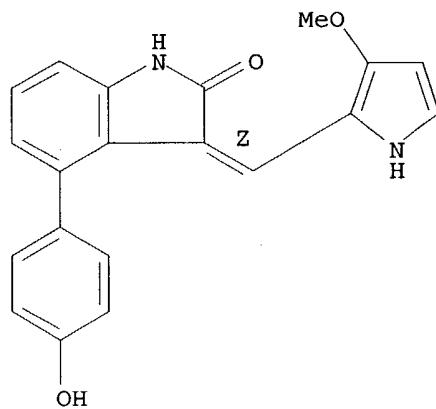
RN 276251-35-5 CAPLUS  
 CN 2H-Indol-2-one, 1,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-4-phenyl-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



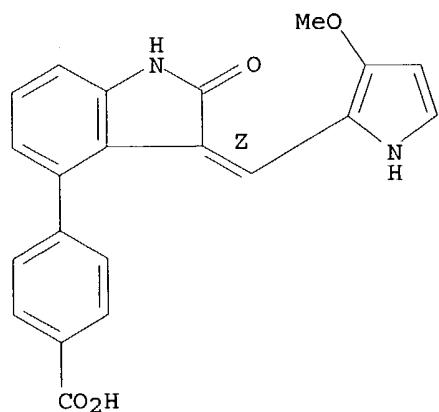
RN 276251-36-6 CAPLUS  
 CN 2H-Indol-2-one, 1,3-dihydro-4-(4-hydroxyphenyl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 276251-37-7 CAPLUS  
 CN Benzoic acid, 4-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-1H-indol-4-yl]- (9CI) (CA INDEX NAME)

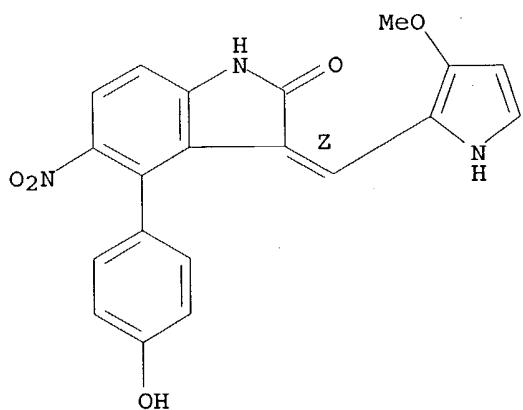
Double bond geometry as shown.



RN 276251-38-8 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(4-hydroxyphenyl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-5-nitro-, (3Z)- (9CI) (CA INDEX NAME)

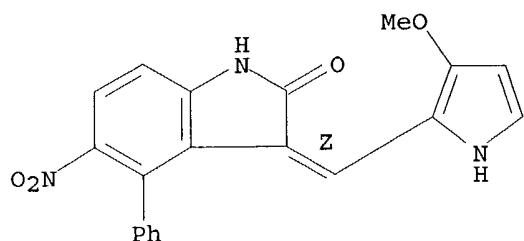
Double bond geometry as shown.



RN 276251-39-9 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-5-nitro-4-phenyl-, (3Z)- (9CI) (CA INDEX NAME)

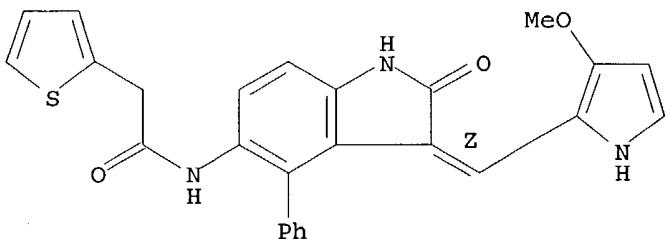
Double bond geometry as shown.



RN 276251-40-2 CAPLUS

CN 2-Thiopheneacetamide, N-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-4-phenyl-1H-indol-5-yl]- (9CI) (CA INDEX NAME)

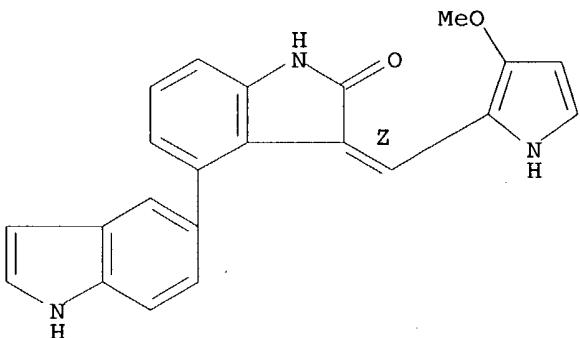
Double bond geometry as shown.



RN 276251-41-3 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(1H-indol-5-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

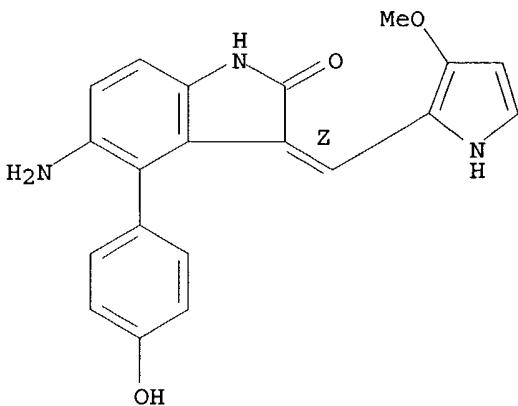
Double bond geometry as shown.



RN 276251-42-4 CAPLUS

CN 2H-Indol-2-one, 5-amino-1,3-dihydro-4-(4-hydroxyphenyl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

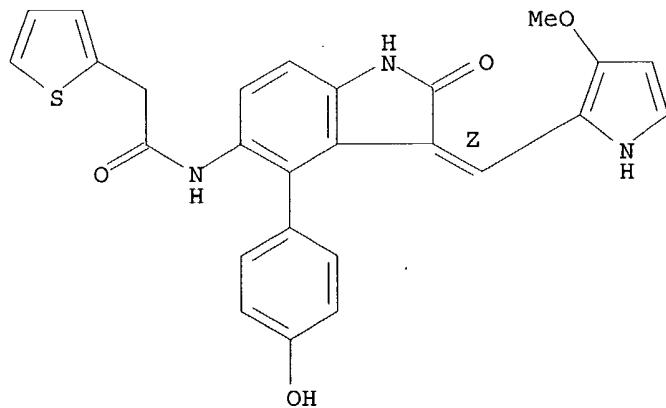
Double bond geometry as shown.



RN 276251-43-5 CAPLUS

CN 2-Thiopheneacetamide, N-[(3Z)-2,3-dihydro-4-(4-hydroxyphenyl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-1H-indol-5-yl]- (9CI) (CA INDEX NAME)

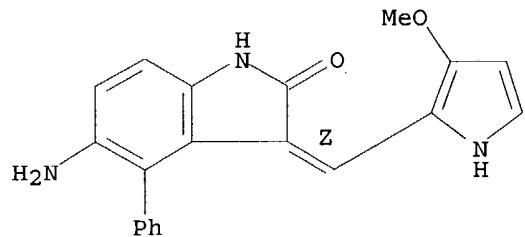
Double bond geometry as shown.



RN 276251-44-6 CAPLUS

CN 2H-Indol-2-one, 5-amino-1,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-4-phenyl-, (3Z)- (9CI) (CA INDEX NAME)

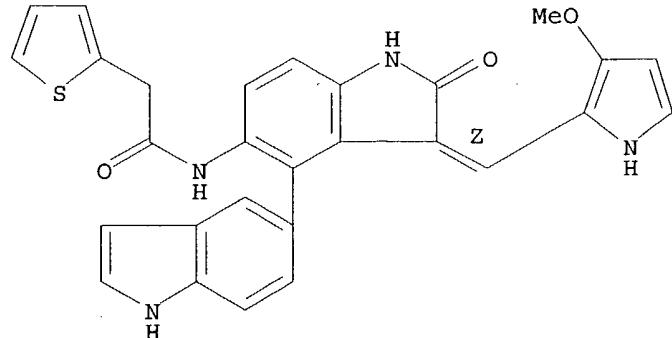
Double bond geometry as shown.



RN 276251-45-7 CAPLUS

CN 2-Thiopheneacetamide, N-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo[4,5'-bi-1H-indol]-5-yl]- (9CI) (CA INDEX NAME)

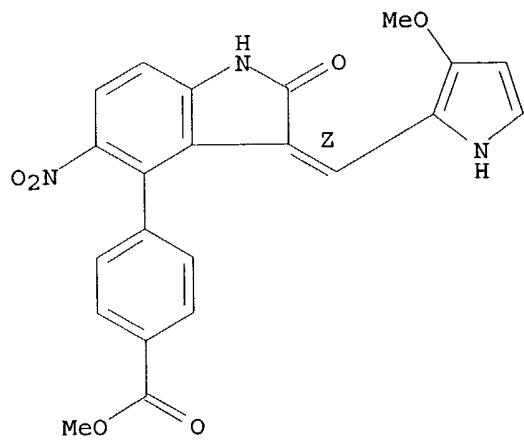
Double bond geometry as shown.



RN 276251-46-8 CAPLUS

CN Benzoic acid, 4-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-5-nitro-2-oxo-1H-indol-4-yl]-, methyl ester (9CI) (CA INDEX NAME)

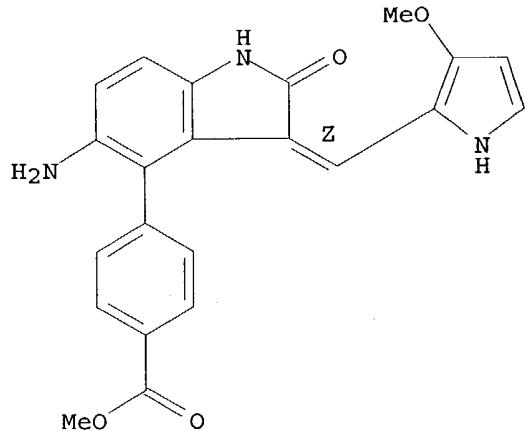
Double bond geometry as shown.



RN 276251-47-9 CAPLUS

CN Benzoic acid, 4-[(3Z)-5-amino-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-1H-indol-4-yl]-, methyl ester (9CI) (CA INDEX NAME)

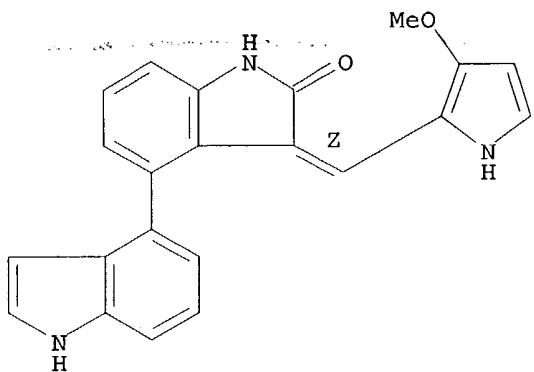
Double bond geometry as shown.



RN 276251-48-0 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(1H-indol-4-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

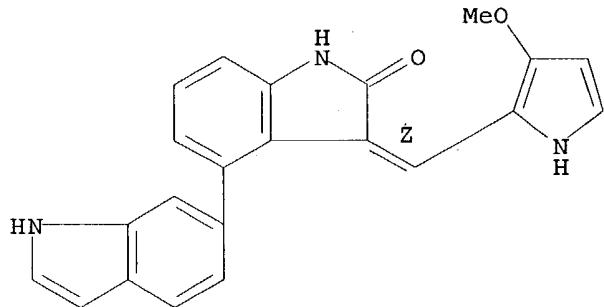
Double bond geometry as shown.



RN 276251-49-1 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(1H-indol-6-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

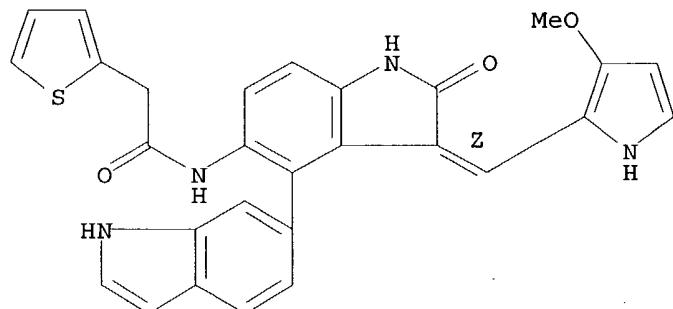
Double bond geometry as shown.



RN 276251-50-4 CAPLUS

CN 2-Thiopheneacetamide, N-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo[4,6'-bi-1H-indol]-5-yl]- (9CI) (CA INDEX NAME)

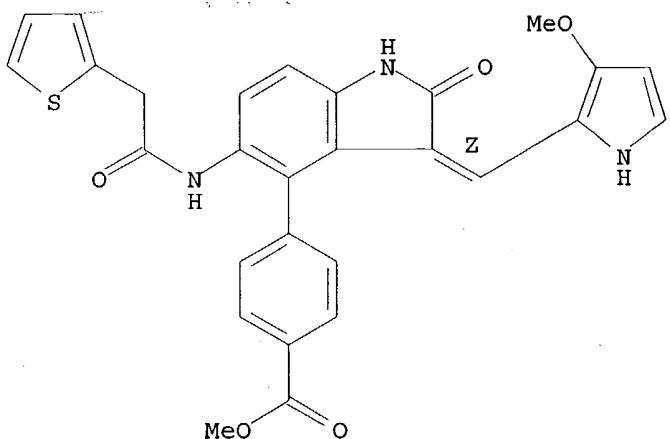
Double bond geometry as shown.



RN 276251-51-5 CAPLUS

CN Benzoic acid, 4-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-5-[(2-thienylacetyl)amino]-1H-indol-4-yl]-, methyl ester (9CI) (CA INDEX NAME)

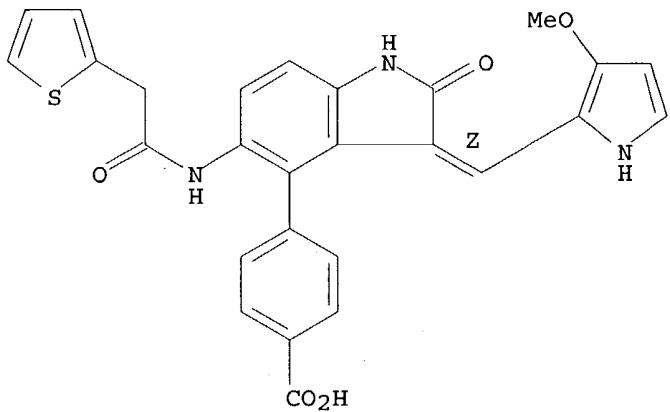
Double bond geometry as shown.



RN 276251-52-6 CAPLUS

CN Benzoic acid, 4-[(3Z)-2,3-dihydro-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-5-[(2-thienylacetyl)amino]-1H-indol-4-yl]- (9CI) (CA INDEX NAME)

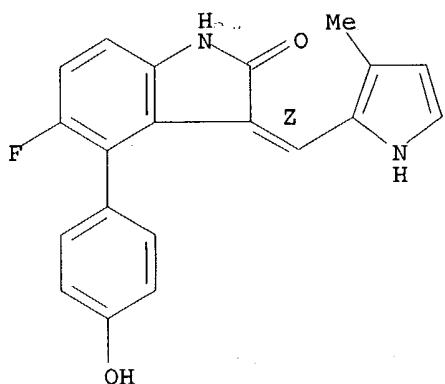
Double bond geometry as shown.



RN 276251-67-3 CAPLUS

CN 2H-Indol-2-one, 5-fluoro-1,3-dihydro-4-(4-hydroxyphenyl)-3-[(3-methyl-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

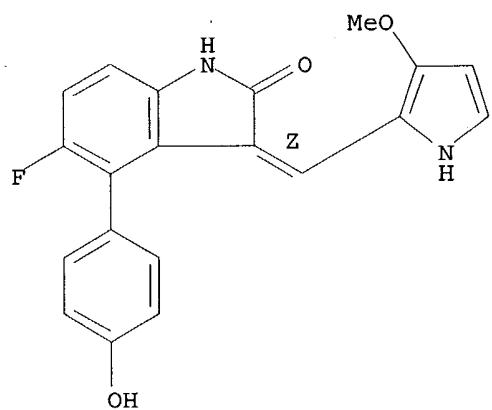
Double bond geometry as shown.



RN 276251-68-4 CAPLUS

CN 2H-Indol-2-one, 5-fluoro-1,3-dihydro-4- (4-hydroxyphenyl) -3- [(3-methoxy-1H-pyrrol-2-yl)methylene] -, (3Z) - (9CI) (CA INDEX NAME)

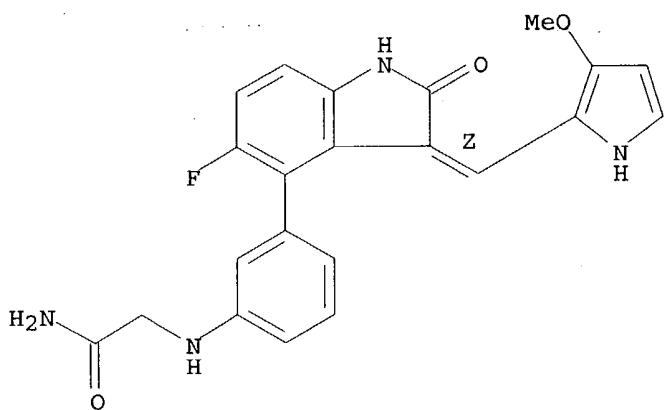
Double bond geometry as shown.



RN 276251-69-5 CAPLUS

CN Acetamide, 2- [ [3- [(3Z)-5-fluoro-2,3-dihydro-3- [(3-methoxy-1H-pyrrol-2-yl)methylene]-2-oxo-1H-indol-4-yl]phenyl]amino] - (9CI) (CA INDEX NAME)

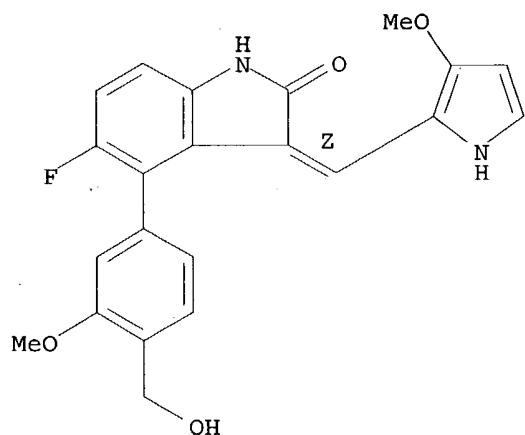
Double bond geometry as shown.



RN 276251-70-8 CAPLUS

CN 2H-Indol-2-one, 5-fluoro-1,3-dihydro-4-[4-(hydroxymethyl)-3-methoxyphenyl]-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

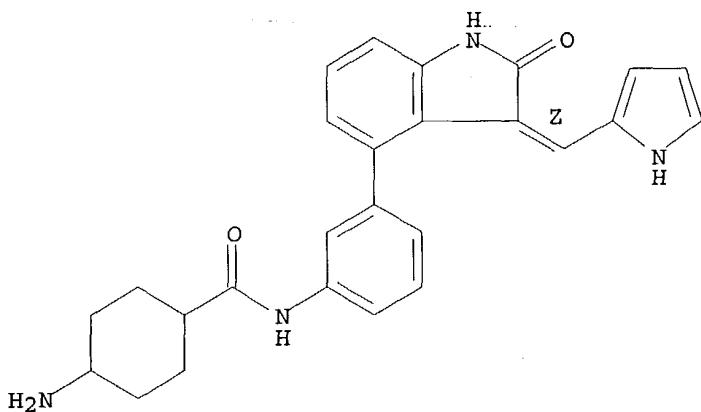
Double bond geometry as shown.



RN 276256-00-9 CAPLUS

CN Cyclohexanecarboxamide, 4-amino-N-[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



IT 276251-76-4P 276251-79-7P 276251-80-0P  
 276251-81-1P 276251-82-2P 276256-01-0P

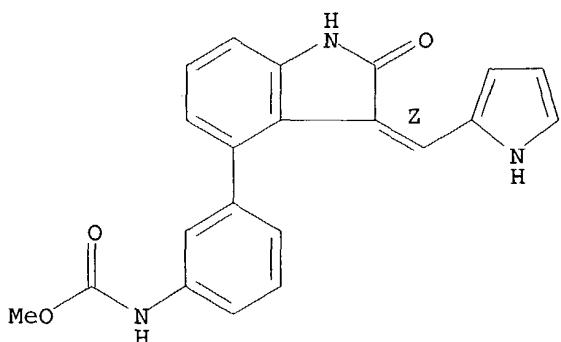
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation of 4-aryl-3-(azolylmethylidene)-2-oxindoles as inhibitors of JNK protein kinases)

RN 276251-76-4 CAPLUS

CN Carbamic acid, [3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

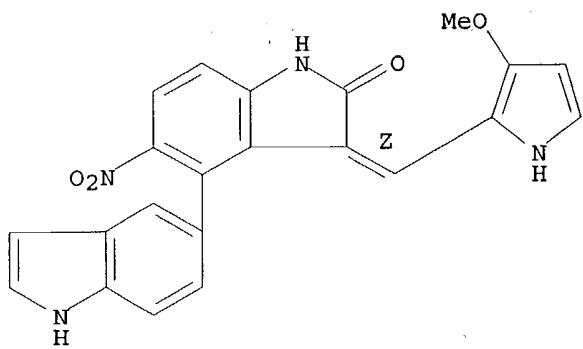
Double bond geometry as shown.



RN 276251-79-7 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(1H-indol-5-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-5-nitro-, (3Z)- (9CI) (CA INDEX NAME)

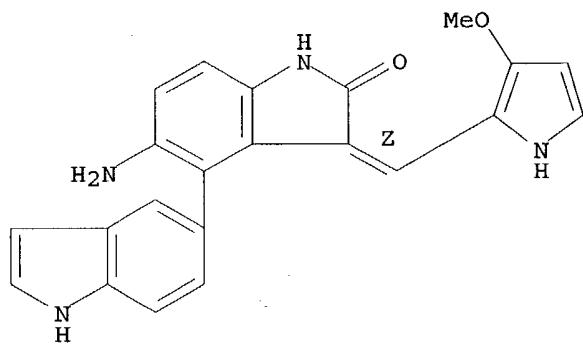
Double bond geometry as shown.



RN 276251-80-0 CAPLUS

CN 2H-Indol-2-one, 5-amino-1,3-dihydro-4-(1H-indol-5-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

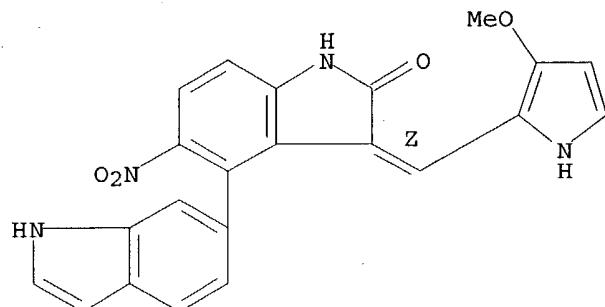
Double bond geometry as shown.



RN 276251-81-1 CAPLUS

CN 2H-Indol-2-one, 1,3-dihydro-4-(1H-indol-6-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-5-nitro-, (3Z)- (9CI) (CA INDEX NAME)

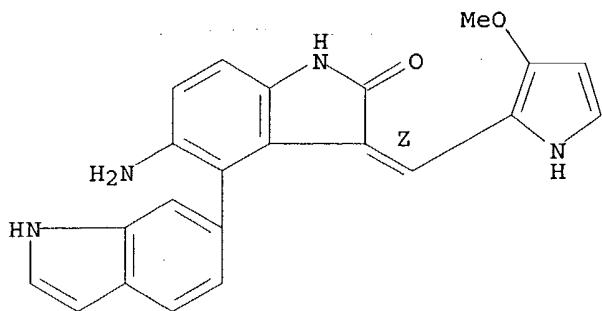
Double bond geometry as shown.



RN 276251-82-2 CAPLUS

CN 2H-Indol-2-one, 5-amino-1,3-dihydro-4-(1H-indol-6-yl)-3-[(3-methoxy-1H-pyrrol-2-yl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

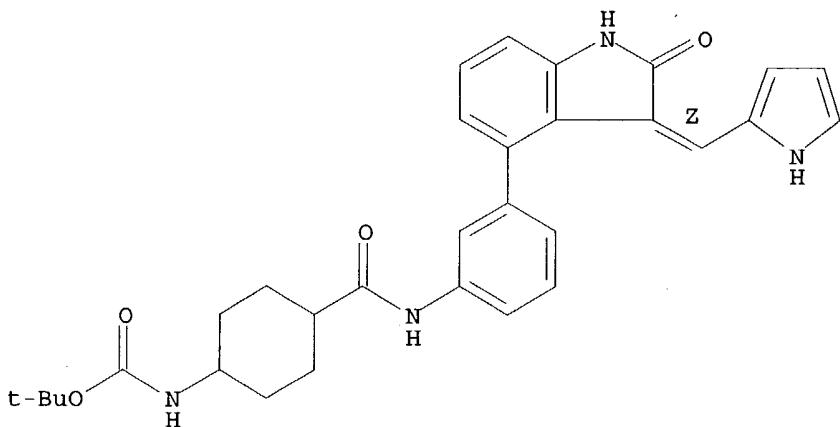
Double bond geometry as shown.



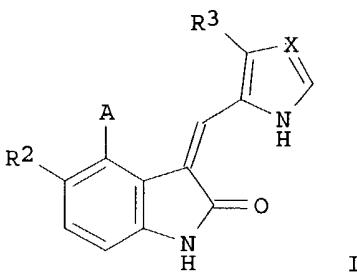
RN 276256-01-0 CAPLUS

CN Carbamic acid, [4-[[3-[(3Z)-2,3-dihydro-2-oxo-3-(1H-pyrrol-2-ylmethylene)-1H-indol-4-yl]phenyl]amino]carbonyl]cyclohexyl]-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

Double bond geometry as shown.



GI



AB Title compds. [I; A = (substituted) aryl, heteroaryl; R2 = H, halo, OR4, NR6R7, COR4, CO2R4, cyano, NO2, SO2R4, SO2NR6R7, etc.; R3 = H, OR4, COR4, CO2R4, CONR6R7, halo, cyano, NR6R7, perfluoroalkyl, (substituted) alkyl,

etc.; R4 = H, (substituted) alkyl, cycloalkyl, heterocyclyl; R6, R7 = H, (substituted) alkyl, cycloalkyl, COR8, CO2R8, SO2R8, etc.; NR6R7 = (substituted) 3-7 membered ring; R8 = H, (substituted) alkyl, aryl, heteroaryl, cycloalkyl; X = N, CH], were prepared. Thus, (Z)-1,3-dihydro-4-iodo-3-[(1H-pyrrol-2-yl)methylene]-2H-indol-2-one (preparation given) was heated with phenylboronic acid, Pd(OAc)2, Et3N, and tri-O-tolylphosphine in DMF at 100° for 24 h to give 85% (Z)-1,3-dihydro-4-phenyl-3-[(1H-pyrrol-2-yl)methylene]-2H-indol-2-one. Tested I inhibited SAPK with IC50<0.15  $\mu$ M.

RE.CNT 2 THERE ARE 2 CITED REFERENCES AVAILABLE FOR THIS RECORD  
ALL CITATIONS AVAILABLE IN THE RE FORMAT